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\*\*\*\*\* Welcome to STN International \*\*\*\*\*

NEWS 1 Web Page for STN Seminar Schedule - N. America  
NEWS 2 AUG 10 Time limit for inactive STN sessions doubles to 40 minutes  
NEWS 3 AUG 18 COMPENDEX indexing changed for the Corporate Source (CS) field  
NEWS 4 AUG 24 ENCOMPLIT/ENCOMPLIT2 reloaded and enhanced  
NEWS 5 AUG 24 CA/CAPLUS enhanced with legal status information for U.S. patents  
NEWS 6 SEP 09 50 Millionth Unique Chemical Substance Recorded in CAS REGISTRY  
NEWS 7 SEP 11 WPIDS, WPINDEX, and WPIX now include Japanese FTERM thesaurus  
NEWS 8 OCT 21 Derwent World Patents Index Coverage of Indian and Taiwanese Content Expanded  
NEWS 9 OCT 21 Derwent World Patents Index enhanced with human translated claims for Chinese Applications and Utility Models  
NEWS 10 NOV 23 Addition of SCAN format to selected STN databases  
NEWS 11 NOV 23 Annual Reload of IFI Databases  
NEWS 12 DEC 01 FRFULL Content and Search Enhancements  
NEWS 13 DEC 01 DGENE, USGENE, and PCTGEN: new percent identity feature for sorting BLAST answer sets  
NEWS 14 DEC 02 Derwent World Patent Index: Japanese FI-TERM thesaurus added  
NEWS 15 DEC 02 PCTGEN enhanced with patent family and legal status display data from INPADOCDB  
NEWS 16 DEC 02 USGENE: Enhanced coverage of bibliographic and sequence information  
NEWS 17 DEC 21 New Indicator Identifies Multiple Basic Patent Records Containing Equivalent Chemical Indexing in CA/CAPLUS  
NEWS 18 JAN 12 Match STN Content and Features to Your Information Needs, Quickly and Conveniently  
NEWS 19 JAN 25 Annual Reload of MEDLINE database  
NEWS 20 FEB 16 STN Express Maintenance Release, Version 8.4.2, Is Now Available for Download  
NEWS 21 FEB 16 Derwent World Patents Index (DWPI) Revises Indexing of Author Abstracts  
NEWS 22 FEB 16 New FASTA Display Formats Added to USGENE and PCTGEN  
NEWS 23 FEB 16 INPADOCDB and INPAFAMDB Enriched with New Content and Features  
NEWS 24 FEB 16 INSPEC Adding Its Own IPC codes and Author's E-mail Addresses

NEWS EXPRESS FEBRUARY 15 10 CURRENT WINDOWS VERSION IS V8.4.2,  
AND CURRENT DISCOVER FILE IS DATED 15 JANUARY 2010.

NEWS HOURS      STN Operating Hours Plus Help Desk Availability  
NEWS LOGIN      Welcome Banner and News Items

Enter NEWS followed by the item number or name to see news on that specific topic.

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\* \* \* \* \* STN Columbus \* \* \* \* \*

FILE 'HOME' ENTERED AT 10:25:44 ON 29 MAR 2010

=> fil reg		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.22	0.22

FILE 'REGISTRY' ENTERED AT 10:26:11 ON 29 MAR 2010  
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 28 MAR 2010 HIGHEST RN 1214990-69-8  
DICTIONARY FILE UPDATES: 28 MAR 2010 HIGHEST RN 1214990-69-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

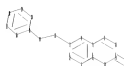
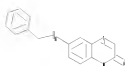
TSCA INFORMATION NOW CURRENT THROUGH January 8, 2010.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

=>  
Uploading C:\Program Files\STNEXP\Queries\10595891latest.str



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chain nodes :
11 13 14
ring nodes :
1 2 3 4 5 6 7 8 9 10 15 16 17 18 19 20
chain bonds :
3-13 9-11 13-14 14-15
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 9-10 15-16 15-20 16-17 17-18
18-19 19-20
exact/norm bonds :
3-13 5-7 6-10 7-8 8-9 9-10 9-11 13-14 14-15
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 15-16 15-20 16-17 17-18 18-19 19-20
isolated ring systems :
containing 1 :

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G1:N,CH

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:CLASS 13:CLASS 14:CLASS 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom

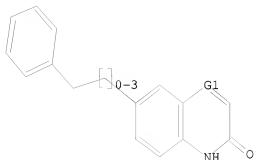
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L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 N,CH

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss full

FULL SEARCH INITIATED 10:26:39 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 79487 TO ITERATE

100.0% PROCESSED 79487 ITERATIONS

291 ANSWERS

SEARCH TIME: 00.00.01

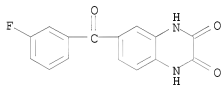
L2 291 SEA SSS FUL L1

=> d scan

L2 291 ANSWERS REGISTRY COPYRIGHT 2010 ACS on STN

IN 2,3-Quinoxalinedione, 6-(3-fluorobenzoyl)-1,4-dihydro-

MF C15 H9 F N2 O3

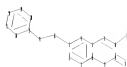
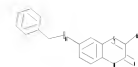


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=>

Uploading C:\Program Files\STNEXP\Queries\10595891last.str



```

chain nodes :
11 13 14 24
ring nodes :
1 2 3 4 5 6 7 8 9 10 15 16 17 18 19 20
chain bonds :
3-13 8-24 9-11 13-14 14-15
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 9-10 15-16 15-20 16-17 17-18
18-19 19-20
exact/norm bonds :
3-13 5-7 8-24 13-14 14-15
exact bonds :
6-10 7-8 8-9 9-10 9-11
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 15-16 15-20 16-17 17-18 18-19 19-20
isolated ring systems :
containing 1 :

```

G1:N,CH

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:CLASS 13:CLASS 14:CLASS 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom
24:CLASS

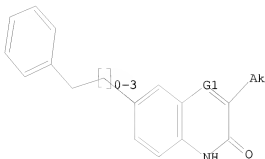
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L3 STRUCTURE UPLOADED

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=> d 13
L3 HAS NO ANSWERS
L3 STR

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G1 N,CH

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SEARCH INITIATED 10:28:45 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 319 TO ITERATE

100.0% PROCESSED 319 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 5309 TO 7451

PROJECTED ANSWERS: 0 TO 0

L4 0 SEA SSS SAM L3

=> s l3 sss full

FULL SEARCH INITIATED 10:28:49 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 6242 TO ITERATE

100.0% PROCESSED 6242 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.02

L5 0 SEA SSS FUL L3

=> fil cap

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

385.04

385.26

FILE 'CAPLUS' ENTERED AT 10:29:32 ON 29 MAR 2010

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FILE COVERS 1907 - 29 Mar 2010 VOL 152 ISS 14  
FILE LAST UPDATED: 28 Mar 2010 (20100328/ED)  
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Dec 2009  
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Dec 2009

CAPLUS now includes complete International Patent Classification (IPC)  
reclassification data for the first quarter of 2010.

CAS Information Use Policies apply and are available at:

<http://www.cas.org/legal/infopolicy.html>

This file contains CAS Registry Numbers for easy and accurate  
substance identification.

=> s l2

L6 18 L2

=> s l2 and (pry<2004 or py<2004)

18 L2  
4301790 PRY<2004  
24050550 PY<2004

L7 15 L2 AND (PRY<2004 OR PY<2004)

=> d l-15 ibib abs hitstr

L7 ANSWER 1 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

ACCESSION NUMBER: 2005:523430 CAPLUS

DOCUMENT NUMBER: 143:60003

TITLE: Preparation of 6-substituted 2-quinolinones and  
2-quinoxalinones as poly(ADP-ribose) polymerase  
inhibitors

INVENTOR(S): Mabire, Dominique Jean-Pierre; Guillemont, Jerome  
Emile Georges; Van Dun, Jacobus Alphonsus Josephus;  
Somers, Maria Victorina Francisca; Wouters, Walter  
Boudewijn Leopold

PATENT ASSIGNEE(S): Janssen Pharmaceutica N. V., Belg.

SOURCE: PCT Int. Appl., 48 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

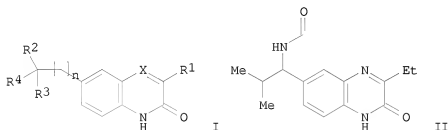
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005054210	A1	20050616	WO 2004-EP13164	20041118 <--
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, GR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
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CA 2546657	A1	20050616	CA 2004-2546657	20041118 <--

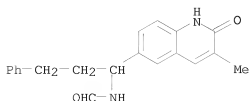
EP 1709012 A1 20061011 EP 2004-819602 20041118 <--  
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK,  
HR, IS, YU  
CN 1890224 A 20070103 CN 2004-80035857 20041118 <--  
BR 2004016532 A 20070109 BR 2004-16532 20041118 <--  
JP 2007513101 T 20070524 JP 2006-541830 20041118 <--  
SG 151249 A1 20090430 SG 2009-1531 20041118 <--  
IN 2006DN03071 A 20070810 IN 2006-DN3071 20060529 <--  
US 20070129375 A1 20070607 US 2006-596086 20060530 <--  
MX 2006006255 A 20060809 MX 2006-6255 20060602 <--  
KR 2006118534 A 20061123 KR 2006-711234 20060608 <--  
NO 2006003028 A 20060628 NO 2006-3028 20060628 <--  
PRIORITY APPLN. INFO.: EP 2003-78859 A 20031205 <--  
WO 2004-EP13164 W 20041118  
ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT  
OTHER SOURCE(S): CASREACT 143:60003; MARPAT 143:60003  
GI



AB The title compds. I [n = 0-2; X = N, CR5; R5 = H or taken together with R1 may form CH:CHCH:CH; R1 = alkyl, thienyl; R2 = H, OH, or taken together with R3 or R4 may form O; R3 = OH, OR8, SR9, etc.; R8 = alkyl, alkylcarbonyl, dialkylaminoalkyl; R9 = dialkylaminoalkyl; R4 = H, alkyl, furanyl, etc.; with the provision], useful for the treatment of a PARP mediated disorder, were prepared E.g., a multi-step synthesis of II, starting from 1-(4-amino-3-nitrophenyl)-2-methyl-1-propanone, was given. The exemplified compds. I were tested in an in vitro assay based on SPA technol. and in an in vitro filtration assay assessing PARP-1 activity (data given). The pharmaceutical composition comprising the compound I is disclosed.

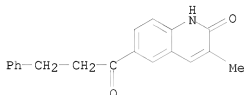
IT 854523-82-3P  
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of 6-substituted 2-quinolinones and 2-quinoxalinones as poly(ADP-ribose) polymerase inhibitors)

RN 854523-82-3 CAPLUS  
CN Formamide, N-[1-(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)-3-phenylpropyl]-  
(CA INDEX NAME)





IT 854524-08-6P  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT  
 (Reactant or reagent)  
 (preparation of 6-substituted 2-quinolinones and 2-quinoxalinones as  
 poly(ADP-ribose) polymerase inhibitors)  
 RN 854524-08-6 CAPLUS  
 CN 2(1H)-Quinolinone, 3-methyl-6-(1-oxo-3-phenylpropyl)- (CA INDEX NAME)



OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD  
 (2 CITINGS)  
 REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

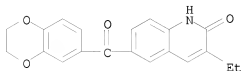
L7 ANSWER 2 OF 15 CAPLUS COPYRIGHT 2010 ACS ON STN  
 ACCESSION NUMBER: 2005:523424 CAPLUS  
 DOCUMENT NUMBER: 143:60001  
 TITLE: Preparation of 6-alkenyl and 6-phenylalkyl substituted  
 2-quinolinones and 2-quinoxalinones as  
 poly(ADP-ribose) polymerase inhibitors  
 INVENTOR(S): Mabire, Dominique Jean-pierre; Guillemont, Jerome  
 Emile Georges; Van Dun, Jacobus Alphonsus Josephus;  
 Somers, Maria Victorina Francisca; Wouters, Walter  
 Boudewijn Leopold  
 PATENT ASSIGNEE(S): Janssen Pharmaceutica N. V., Belg.  
 SOURCE: PCI Int. Appl., 102 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

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WO 2005054201	A1	20050616	WO 2004-EP13163	20041118 <--
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RW:	BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
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EP 1687277	A1	20060809	EP 2004-819601	20041118 <--
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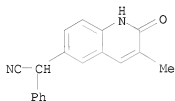
RN 854532-59-5 CAPLUS

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(CA INDEX NAME)



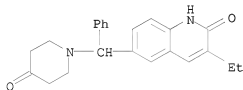
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CN 6-Quinolineacetonitrile, 1,2-dihydro-3-methyl-2-oxo- $\alpha$ -phenyl- (CA  
INDEX NAME)



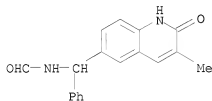
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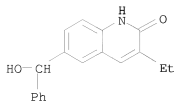
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CN Formamide, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)phenylmethyl]- (CA  
INDEX NAME)



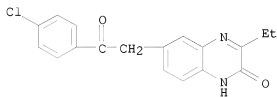
RN 854534-00-2 CAPLUS

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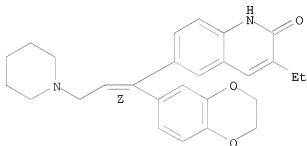
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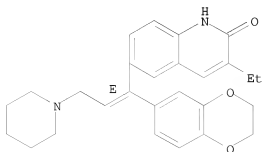
Double bond geometry as shown.



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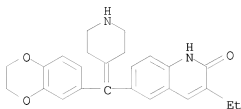
CN 2(1H)-Quinolone, 6-[(1E)-1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(1-piperidinyl)-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

Double bond geometry as shown.



RN 854534-19-3 CAPLUS

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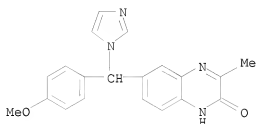
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	854532-64-2P	854532-65-3P	854532-66-4P
	854532-67-5P	854532-69-7P	854532-70-0P
	854532-71-1P	854532-72-2P	854532-73-3P
	854532-74-4P	854532-75-5P	854532-76-6P
	854532-77-7P	854532-78-8P	854532-79-9P
	854532-80-2P	854532-81-3P	854532-82-4P
	854532-83-5P	854532-84-6P	854532-85-7P
	854532-86-8P	854532-87-9P	854532-89-1P
	854532-92-6P	854532-93-7P	854532-94-8P
	854532-95-9P	854532-96-0P	854532-97-1P
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	854533-09-8P	854533-14-5P	854533-16-7P
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	854533-38-3P	854533-39-4P	854533-40-7P
	854533-41-8P	854533-42-9P	854533-43-0P
	854533-44-1P	854533-45-2P	854533-46-3P
	854533-47-4P	854533-48-5P	854533-49-6P
	854533-51-0P	854533-53-2P	854533-54-3P
	854533-55-4P	854533-56-5P	854533-57-6P
	854533-58-7P	854533-59-8P	854533-60-1P
	854533-62-3P	854533-65-6P	854533-67-8P
	854533-69-0P	854533-71-4P	854533-73-6P
	854533-75-8P	854533-79-2P	854533-81-6P
	854533-83-8P	854533-85-0P	854533-87-2P
	854533-89-4P	854533-91-8P	854533-93-0P
	854533-97-4P	854533-98-5P	854533-99-6P
	854534-01-3P	854534-02-4P	854534-04-6P
	854534-05-7P	854534-06-8P	854534-07-9P
	854534-08-0P	854534-09-1P	854534-10-4P
	854534-11-5P	854534-12-6P	854534-13-7P
	854534-14-8P	854534-15-9P	854534-16-0P
	854534-20-6P	854534-21-7P	854534-22-8P
	854534-23-9P	854534-24-0P	854534-25-1P
	854534-26-2P	854534-27-3P	854534-28-4P
	854534-29-5P	854534-30-8P	854534-31-9P
	854534-32-0P	854534-33-1P	854535-35-6P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of 6-alkenyl and 6-phenylalkyl substituted 2-quinolinones and 2-quinoxalinones as poly(ADP-ribose) polymerase inhibitors)

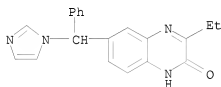
RN 130346-68-8 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[1H-imidazol-1-yl(4-methoxyphenyl)methyl]-3-methyl-  
(CA INDEX NAME)



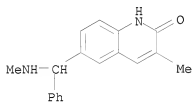
RN 130346-69-9 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-(1H-imidazol-1-ylphenylmethyl)- (CA INDEX  
NAME)



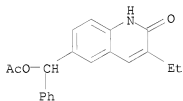
RN 854532-58-4 CAPLUS

CN 2(1H)-Quinolinone, 3-methyl-6-[(methylamino)phenylmethyl]- (CA INDEX  
NAME)



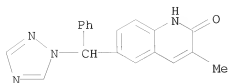
RN 854532-60-8 CAPLUS

CN 2(1H)-Quinolinone, 6-[(acetyloxy)phenylmethyl]-3-ethyl- (CA INDEX NAME)



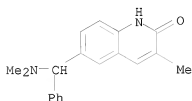
RN 854532-62-0 CAPLUS

CN 2(1H)-Quinolinone, 3-methyl-6-(phenyl-1H-1,2,4-triazol-1-ylmethyl)- (CA  
INDEX NAME)



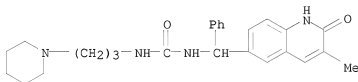
RN 854532-63-1 CAPLUS

CN 2(1H)-Quinolinsonone, 6-[(dimethylamino)phenylmethyl]-3-methyl- (CA INDEX NAME)



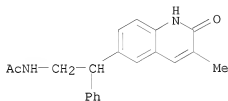
RN 854532-64-2 CAPLUS

CN Urea, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)phenylmethyl]-N'-[3-(1-piperidinyl)propyl]- (CA INDEX NAME)



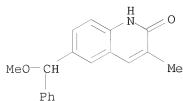
RN 854532-65-3 CAPLUS

CN Acetamide, N-[2-(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)-2-phenylethyl]- (CA INDEX NAME)

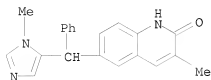


RN 854532-66-4 CAPLUS

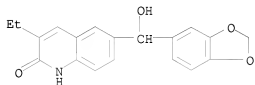
CN 2(1H)-Quinolinsonone, 6-(methoxyphenylmethyl)-3-methyl- (CA INDEX NAME)



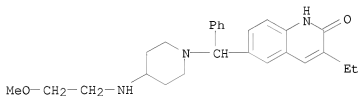
RN 854532-67-5 CAPLUS  
 CN 2(1H)-Quinolinone, 3-methyl-6-[(1-methyl-1H-imidazol-5-yl)phenylmethyl]-  
 (CA INDEX NAME)



RN 854532-69-7 CAPLUS  
 CN 2(1H)-Quinolinone, 6-(1,3-benzodioxol-5-ylhydroxymethyl)-3-ethyl- (CA  
 INDEX NAME)



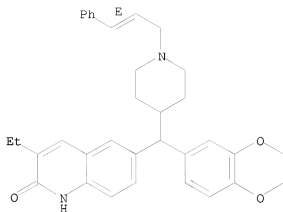
RN 854532-70-0 CAPLUS  
 CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-  
 piperidinyl]phenylmethyl]- (CA INDEX NAME)



RN 854532-71-1 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[1-[(2E)-3-phenyl-  
 2-propen-1-yl]-4-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)

Double bond geometry as shown.

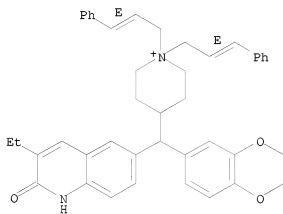




RN 854532-72-2 CAPLUS

CN Piperidinium, 4-[(2,3-dihydro-1,4-benzodioxin-6-yl)(3-ethyl-1,2-dihydro-2-oxo-6-quinolinyl)methyl]-1,1-bis[(2E)-3-phenyl-2-propen-1-yl]-, chloride (1:1) (CA INDEX NAME)

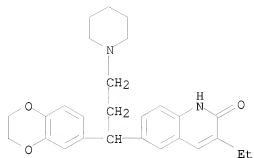
Double bond geometry as shown.



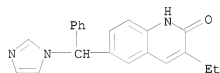
● Cl<sup>-</sup>

RN 854532-73-3 CAPLUS

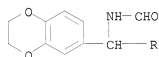
CN 2(1H)-Quinolinone, 6-[1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(1-piperidinyl)propyl]-3-ethyl- (CA INDEX NAME)



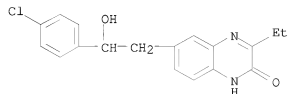
RN 854532-74-4 CAPLUS  
 CN 2(1H)-Quinolinone, 3-ethyl-6-(1H-imidazol-1-ylphenylmethyl)- (CA INDEX NAME)



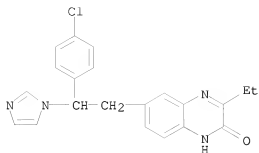
RN 854532-75-5 CAPLUS  
 CN Formamide, N-[2,3-dihydro-1,4-benzodioxin-6-yl] (3-ethyl-1,2-dihydro-2-oxo-6-quinolinyl)methyl)- (CA INDEX NAME)



RN 854532-76-6 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-hydroxyethyl]-3-ethyl- (CA INDEX NAME)

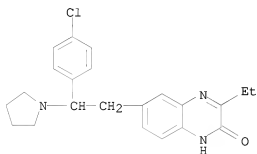


RN 854532-77-7 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-(1H-imidazol-1-yl)ethyl]-3-ethyl- (CA INDEX NAME)



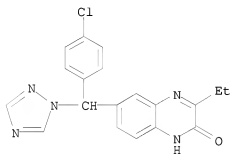
RN 854532-78-8 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-(1-pyrrolidinyl)ethyl]-3-ethyl- (CA INDEX NAME)



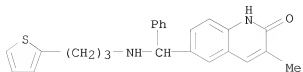
RN 854532-79-9 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-1,2,4-triazol-1-ylmethyl]-3-ethyl- (CA INDEX NAME)



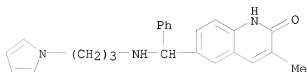
RN 854532-80-2 CAPLUS

CN 2(1H)-Quinololinone, 3-methyl-6-[phenyl[[3-(2-thienyl)propyl]amino]methyl]- (CA INDEX NAME)



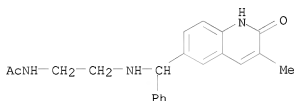
RN 854532-81-3 CAPLUS

CN 2(1H)-Quinolinone, 3-methyl-6-[phenyl[[3-(1H-pyrrol-1-yl)propyl]amino]methyl]- (CA INDEX NAME)



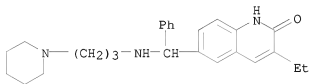
RN 854532-82-4 CAPLUS

CN Acetamide, N-[2-[[[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)phenylmethyl]amino]ethyl]- (CA INDEX NAME)



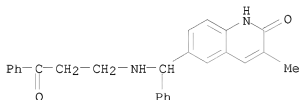
RN 854532-83-5 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[phenyl[[3-(1-piperidinyl)propyl]amino]methyl]- (CA INDEX NAME)



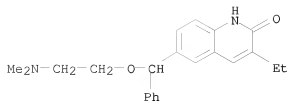
RN 854532-84-6 CAPLUS

CN 2(1H)-Quinolinone, 3-methyl-6-[[[(3-oxo-3-phenylpropyl)amino]phenylmethyl]- (CA INDEX NAME)



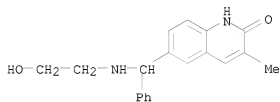
RN 854532-85-7 CAPLUS

CN 2(1H)-Quinolinone, 6-[[[2-(dimethylamino)ethoxy]phenylmethyl]-3-ethyl- (CA INDEX NAME)



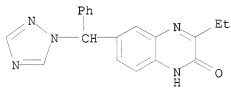
RN 854532-86-8 CAPLUS

CN 2(1H)-Quinolinone, 6-[[[(2-hydroxyethyl)amino]phenylmethyl]-3-methyl- (CA  
INDEX NAME)



RN 854532-87-9 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-(phenyl-1H-1,2,4-triazol-1-ylmethyl)- (CA  
INDEX NAME)



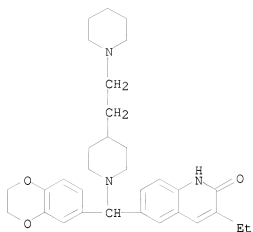
RN 854532-89-1 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[4-[2-(1-piperidinyl)ethyl]-1-piperidinyl)methyl]-3-ethyl-, ethanedioate (2:5) (CA  
INDEX NAME)

CM 1

CRN 854532-88-0

CMF C32 H41 N3 O3



CM 2

CRN 144-62-7

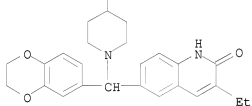
CMF C2 H2 O4



RN 854532-92-6 CAPLUS

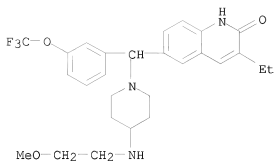
CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)

MeO-CH<sub>2</sub>-CH<sub>2</sub>-NH



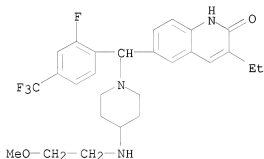
RN 854532-93-7 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidinyl][3-(trifluoromethoxy)phenyl)methyl]- (CA INDEX NAME)



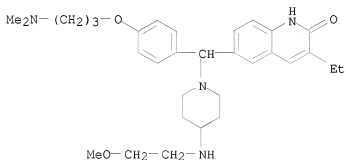
RN 854532-94-8 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[[2-fluoro-4-(trifluoromethyl)phenyl][4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]- (CA INDEX NAME)



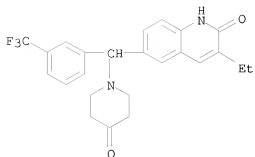
RN 854532-95-9 CAPLUS

CN 2(1H)-Quinolinone, 6-[[4-[3-(dimethylamino)propoxy]phenyl][4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



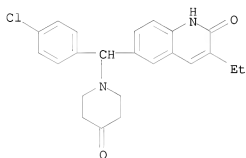
RN 854532-96-0 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(4-oxo-1-piperidinyl)[3-(trifluoromethyl)phenyl)methyl]- (CA INDEX NAME)



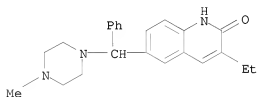
RN 854532-97-1 CAPLUS

CN 2(1H)-Quinolinone, 6-[(4-chlorophenyl)(4-oxo-1-piperidinyl)methyl]-3-ethyl-  
(CA INDEX NAME)



RN 854532-98-2 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(4-methyl-1-piperazinyl)phenylmethyl]-  
(CA INDEX NAME)



RN 854533-00-9 CAPLUS

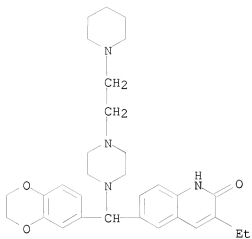
CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[4-[2-(1-piperidinyl)ethyl]-1-piperazinyl)methyl]-3-ethyl-, ethanedioate (1:3)  
(CA INDEX NAME)

CM 1

CRN 854532-99-3

CMF C31 H40 N4 O3





CM 2

CRN 144-62-7

CMF C2 H2 O4



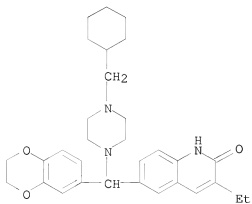
RN 854533-02-1 CAPLUS

CN 2-(1H)-Quinolinone, 6-[[4-(cyclohexylmethyl)-1-piperazinyl](2,3-dihydro-1,4-benzodioxin-6-yl)methyl]-3-ethyl-, ethanedioate (2:3) (CA INDEX NAME)

CM 1

CRN 854533-01-0

CMF C31 H39 N3 O3



CM 2

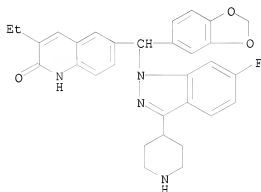
CRN 144-62-7

CMF C2 H2 O4



RN 854533-04-3 CAPLUS

CN 2(1H)-Quinolinone, 6-[(1,3-benzodioxol-5-yl[6-fluoro-3-(4-piperidinyl)-1H-indazol-1-yl]methyl]-3-ethyl- (CA INDEX NAME)



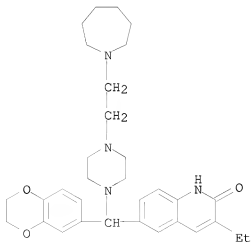
RN 854533-06-5 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[4-[2-(hexahydro-1H-azepin-1-yl)ethyl]-1-piperazinyl]methyl]-3-ethyl-, ethanedioate (2:5) (CA INDEX NAME)

CM 1

CRN 854533-05-4

CMF C32 H42 N4 O3



CM 2

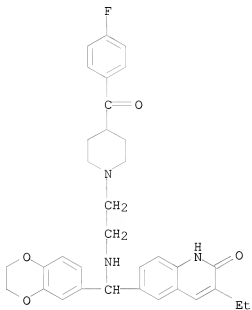
CRN 144-62-7

CMF C2 H2 O4



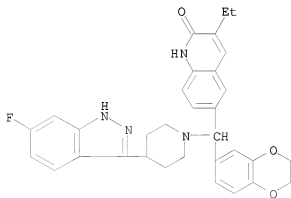
RN 854533-07-6 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[[2-[4-(4-fluorobenzoyl)-1-piperidinyl]ethyl]amino]methyl]-3-ethyl- (CA INDEX NAME)



RN 854533-09-8 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[4-(6-fluoro-1H-indazol-3-yl)-1-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)



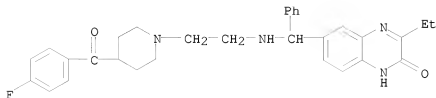
RN 854533-14-5 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-[[[2-[4-(4-fluorobenzoyl)-1-piperidinyl]ethyl]amino]phenylmethyl]-, ethanedioate (2:5) (CA INDEX NAME)

CM 1

CRN 854533-13-4

CMF C31 H33 F N4 O2



CM 2

CRN 144-62-7

CMF C2 H2 O4



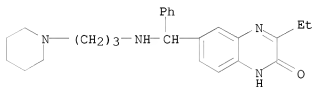
RN 854533-16-7 CAPLUS

CN 2-(1H)-Quinoxalinone, 3-ethyl-6-[phenyl[[3-(1-piperidinyl)propyl]amino]methyl]-, ethanedioate (2:5) (CA INDEX NAME)

CM 1

CRN 854533-15-6

CMF C25 H32 N4 O



CM 2

CRN 144-62-7

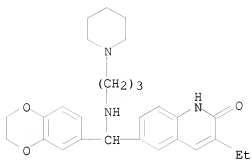
CMF C2 H2 O4



RN 854533-18-9 CAPLUS

CN 2-(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[[3-(1-piperidinyl)propyl]amino]methyl]-3-ethyl-, ethanedioate (1:2) (CA INDEX NAME)

CM 1  
 CRN 854533-17-8  
 CMF C28 H35 N3 O3

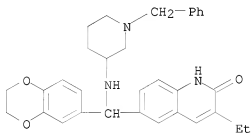


CM 2  
 CRN 144-62-7  
 CMF C2 H2 O4



RN 854533-20-3 CAPLUS  
 CN 2-(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[[1-(phenylmethyl)-3-piperidinylamino]methyl]-3-ethyl-, ethanedioate (2:5) (CA INDEX NAME)

CM 1  
 CRN 854533-19-0  
 CMF C32 H35 N3 O3

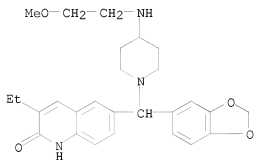


CM 2  
 CRN 144-62-7  
 CMF C2 H2 O4



RN 854533-21-4 CAPLUS

CN 2(1H)-Quinololinone, 6-[1,3-benzodioxol-5-yl[4-[(2-methoxyethyl)amino]-1-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)



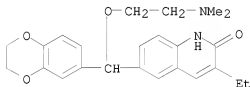
RN 854533-23-6 CAPLUS

CN 2(1H)-Quinololinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[2-(dimethylamino)ethoxy]methyl]-3-ethyl-, ethanedioate (2:3) (CA INDEX NAME)

CM 1

CRN 854533-22-5

CMF C24 H28 N2 O4



CM 2

CRN 144-62-7

CMF C2 H2 O4

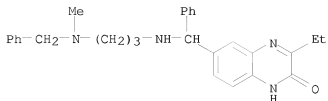


RN 854533-25-8 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-[[[3-methyl(phenylmethyl)amino]propyl]amino]phenylmethyl]-, ethanedioate (1:2) (CA INDEX NAME)

CM 1

CRN 854533-24-7  
CMF C28 H32 N4 O



CM 2

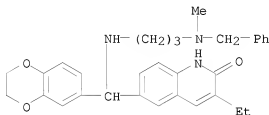
CRN 144-62-7  
CMF C2 H2 O4



RN 854533-27-0 CAPLUS  
CN 2-(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[[3-methyl(phenylmethyl)amino]propyl]amino]methyl-3-ethyl-, ethanedioate (1:2) (CA INDEX NAME)

CM 1

CRN 854533-26-9  
CMF C31 H35 N3 O3



CM 2

CRN 144-62-7  
CMF C2 H2 O4

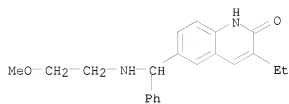


RN 854533-29-2 CAPLUS  
CN 2-(1H)-Quinolinone, 3-ethyl-6-[[2-methoxyethyl]amino]phenylmethyl-, ethanedioate (1:2) (CA INDEX NAME)

CM 1

CRN 854533-28-1

CMF C21 H24 N2 O2



CM 2

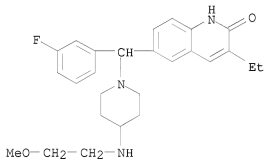
CRN 144-62-7

CMF C2 H2 O4



RN 854533-30-5 CAPLUS

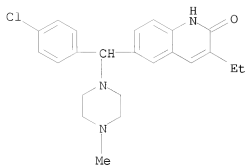
CN 2-(1H)-Quinolinone, 3-ethyl-6-[(3-fluorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]- (CA INDEX NAME)



RN 854533-31-6 CAPLUS

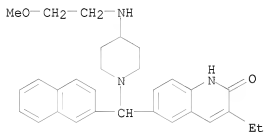
CN 2-(1H)-Quinolinone, 6-[(4-chlorophenyl)(4-methyl-1-piperazinyl)methyl]-3-ethyl- (CA INDEX NAME)





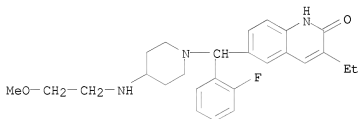
RN 854533-32-7 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidiny]-2-naphthalenylmethyl]- (CA INDEX NAME)



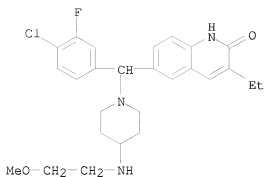
RN 854533-33-8 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(2-fluorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidiny]methyl]- (CA INDEX NAME)



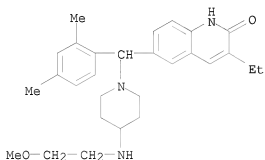
RN 854533-34-9 CAPLUS

CN 2(1H)-Quinolinone, 6-[(4-chloro-3-fluorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidiny]methyl]-3-ethyl- (CA INDEX NAME)



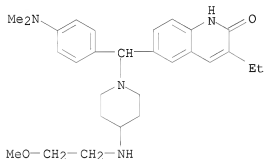
RN 854533-35-0 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,4-dimethylphenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)



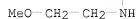
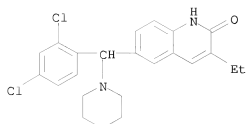
RN 854533-36-1 CAPLUS

CN 2(1H)-Quinolinone, 6-[[4-(dimethylamino)phenyl][4-[(2-methoxyethyl)amino]-1-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)



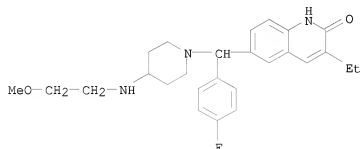
RN 854533-37-2 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,4-dichlorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl]methyl]-3-ethyl- (CA INDEX NAME)



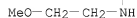
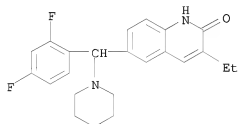
RN 854533-38-3 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(4-fluorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]- (CA INDEX NAME)



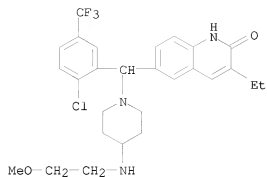
RN 854533-39-4 CAPLUS

CN 2(1H)-Quinolinone, 6-[[2,4-difluorophenyl][4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



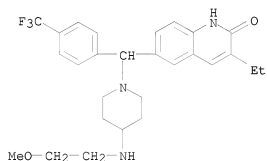
RN 854533-40-7 CAPLUS

CN 2(1H)-Quinolinone, 6-[[2-chloro-5-(trifluoromethyl)phenyl][4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



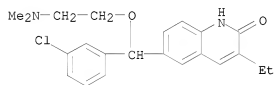
RN 854533-41-8 CAPLUS

CN 2(1H)-Quinolinsonone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidinyl][4-(trifluoromethyl)phenyl)methyl]- (CA INDEX NAME)



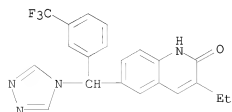
RN 854533-42-9 CAPLUS

CN 2(1H)-Quinolinsonone, 6-[(3-chlorophenyl)[2-(dimethylamino)ethoxy)methyl]-3-ethyl- (CA INDEX NAME)



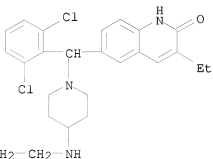
RN 854533-43-0 CAPLUS

CN 2(1H)-Quinolinsonone, 3-ethyl-6-[[4H-1,2,4-triazol-4-yl][3-(trifluoromethyl)phenyl)methyl]- (CA INDEX NAME)



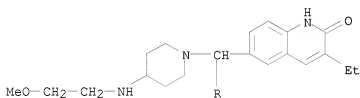
RN 854533-44-1 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2,6-dichlorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



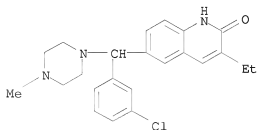
RN 854533-45-2 CAPLUS

CN 2(1H)-Quinolinone, 6-[(2-chlorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



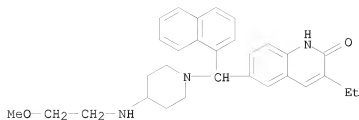
RN 854533-46-3 CAPLUS

CN 2(1H)-Quinolinone, 6-[(3-chlorophenyl)(4-methyl-1-piperazinyl)methyl]-3-ethyl- (CA INDEX NAME)



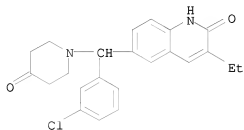
RN 854533-47-4 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidinyl]-1-naphthalenylmethyl]- (CA INDEX NAME)



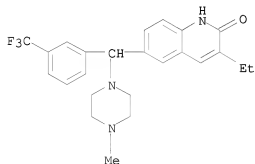
RN 854533-48-5 CAPLUS

CN 2(1H)-Quinolinone, 6-[(3-chlorophenyl)(4-oxo-1-piperidinyl)methyl]-3-ethyl-  
(CA INDEX NAME)



RN 854533-49-6 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(4-methyl-1-piperazinyl)(3-  
(trifluoromethyl)phenyl)methyl]- (CA INDEX NAME)



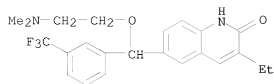
RN 854533-51-0 CAPLUS

CN 2(1H)-Quinolinone, 6-[[2-(dimethylamino)ethoxy](3-  
(trifluoromethyl)phenyl)methyl]-3-ethyl-, ethanedioate (1:1) (CA INDEX  
NAME)

CM 1

CRN 854533-50-9

CMF C23 H25 F3 N2 O2



CM 2

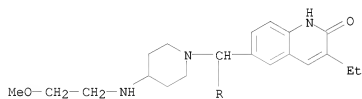
CRN 144-62-7

CMF C2 H2 O4



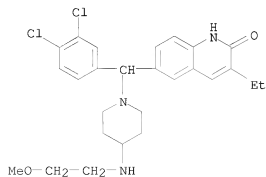
RN 854533-53-2 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidinyl](2-methylphenyl)methyl]- (CA INDEX NAME)

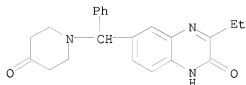


RN 854533-54-3 CAPLUS

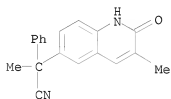
CN 2(1H)-Quinolinone, 6-[(3,4-dichlorophenyl)[4-[(2-methoxyethyl)amino]-1-piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)



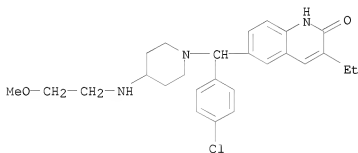
RN 854533-55-4 CAPLUS  
 CN 2(1H)-Quinoxalinone, 3-ethyl-6-[(4-oxo-1-piperidinyl)phenylmethyl]- (CA  
 INDEX NAME)



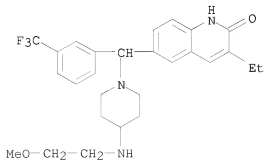
RN 854533-56-5 CAPLUS  
 CN 6-Quinolineacetonitrile, 1,2-dihydro- $\alpha$ ,3-dimethyl-2-oxo- $\alpha$ -  
 phenyl- (CA INDEX NAME)



RN 854533-57-6 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[(4-chlorophenyl)[4-[(2-methoxyethyl)amino]-1-  
 piperidinyl)methyl]-3-ethyl- (CA INDEX NAME)

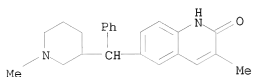


RN 854533-58-7 CAPLUS  
 CN 2(1H)-Quinolinone, 3-ethyl-6-[[4-[(2-methoxyethyl)amino]-1-piperidinyl][3-  
 (trifluoromethyl)phenyl)methyl]- (CA INDEX NAME)

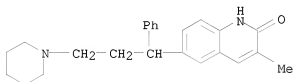




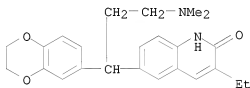
RN 854533-59-8 CAPLUS  
 CN 2(1H)-Quinolinone, 3-methyl-6-[(1-methyl-3-piperidinyl)phenylmethyl]- (CA INDEX NAME)



RN 854533-60-1 CAPLUS  
 CN 2(1H)-Quinolinone, 3-methyl-6-[1-phenyl-3-(1-piperidinyl)propyl]- (CA INDEX NAME)



RN 854533-62-3 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(dimethylamino)propyl]-3-ethyl- (CA INDEX NAME)

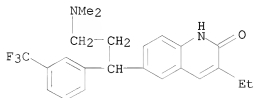


RN 854533-65-6 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[3-(dimethylamino)-1-[3-(trifluoromethyl)phenyl]propyl]-3-ethyl-, ethanedioate (1:1) (CA INDEX NAME)

CM 1

CRN 854533-64-5

CMF C23 H25 F3 N2 O

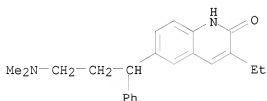


CM 2

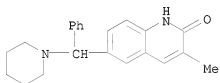
CRN 144-62-7  
CMF C2 H2 O4



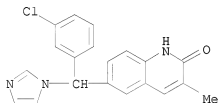
RN 854533-67-8 CAPLUS  
CN 2(1H)-Quinolinone, 6-[3-(dimethylamino)-1-phenylpropyl]-3-ethyl- (CA INDEX NAME)



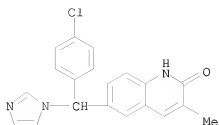
RN 854533-69-0 CAPLUS  
CN 2(1H)-Quinolinone, 3-methyl-6-(phenyl-1-piperidinylmethyl)- (CA INDEX NAME)



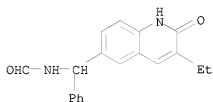
RN 854533-71-4 CAPLUS  
CN 2(1H)-Quinolinone, 6-[(3-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-methyl- (CA INDEX NAME)



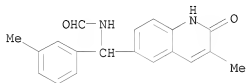
RN 854533-73-6 CAPLUS  
CN 2(1H)-Quinolinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-methyl- (CA INDEX NAME)



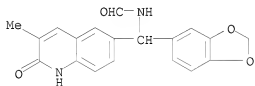
RN 854533-75-8 CAPLUS  
 CN Formamide, N-[(3-ethyl-1,2-dihydro-2-oxo-6-quinolinyl)phenylmethyl]- (CA INDEX NAME)



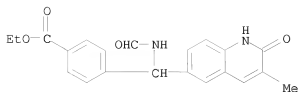
RN 854533-79-2 CAPLUS  
 CN Formamide, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)(3-methylphenyl)methyl]- (CA INDEX NAME)



RN 854533-81-6 CAPLUS  
 CN Formamide, N-[1,3-benzodioxol-5-yl(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)methyl]- (CA INDEX NAME)

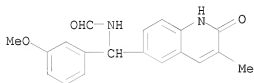


RN 854533-83-8 CAPLUS  
 CN Benzoic acid, 4-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)(formylamino)methyl]-, ethyl ester (CA INDEX NAME)



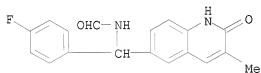
RN 854533-85-0 CAPLUS

CN Formamide, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyne)(3-methoxyphenyl)methyl]- (CA INDEX NAME)



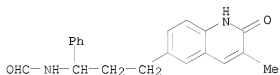
RN 854533-87-2 CAPLUS

CN Formamide, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyne)(4-methoxyphenyl)methyl]- (CA INDEX NAME)



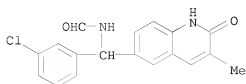
RN 854533-89-4 CAPLUS

CN Formamide, N-[3-(1,2-dihydro-3-methyl-2-oxo-6-quinolinyne)-1-phenylpropyl]- (CA INDEX NAME)



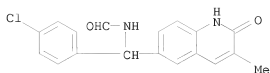
RN 854533-91-8 CAPLUS

CN Formamide, N-[(3-chlorophenyl)(1,2-dihydro-3-methyl-2-oxo-6-quinolinyne)methyl]- (CA INDEX NAME)

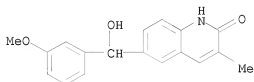


RN 854533-93-0 CAPLUS

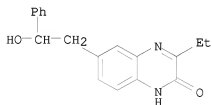
CN Formamide, N-[(4-chlorophenyl)(1,2-dihydro-3-methyl-2-oxo-6-quinolinyne)methyl]- (CA INDEX NAME)



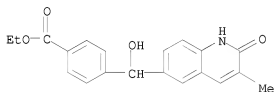
RN 854533-97-4 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[(hydroxy(3-methoxyphenyl)methyl)-3-methyl- (CA INDEX NAME)



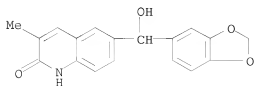
RN 854533-98-5 CAPLUS  
 CN 2(1H)-Quinoxalinone, 3-ethyl-6-(2-hydroxy-2-phenylethyl)- (CA INDEX NAME)



RN 854533-99-6 CAPLUS  
 CN Benzoic acid, 4-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)hydroxymethyl]-, ethyl ester (CA INDEX NAME)

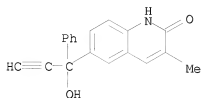


RN 854534-01-3 CAPLUS  
 CN 2(1H)-Quinolinone, 6-[(1,3-benzodioxol-5-yl)hydroxymethyl)-3-methyl- (CA INDEX NAME)



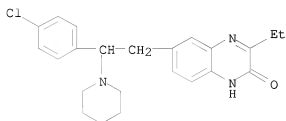
RN 854534-02-4 CAPLUS

CN 2(1H)-Quinolinone, 6-(1-hydroxy-1-phenyl-2-propyn-1-yl)-3-methyl- (CA INDEX NAME)



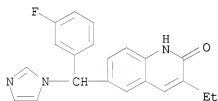
RN 854534-04-6 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-(1-piperidinyl)ethyl]-3-ethyl- (CA INDEX NAME)



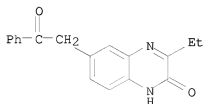
RN 854534-05-7 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[(3-fluorophenyl)-1H-imidazol-1-ylmethyl]- (CA INDEX NAME)



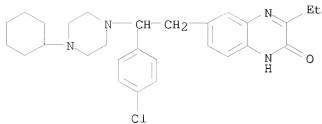
RN 854534-06-8 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-(2-oxo-2-phenylethyl)- (CA INDEX NAME)

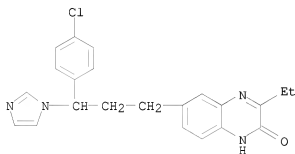


RN 854534-07-9 CAPLUS

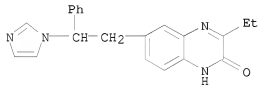
CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-(4-cyclohexyl-1-piperazinyl)ethyl]-3-ethyl- (CA INDEX NAME)



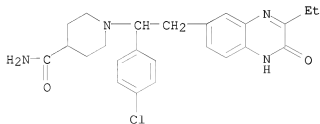
RN 854534-08-0 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[3-(4-chlorophenyl)-3-(1H-imidazol-1-yl)propyl]-3-ethyl- (CA INDEX NAME)



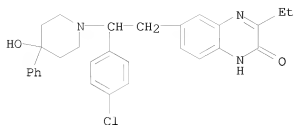
RN 854534-09-1 CAPLUS  
 CN 2(1H)-Quinoxalinone, 3-ethyl-6-[2-(1H-imidazol-1-yl)-2-phenylethyl]- (CA INDEX NAME)



RN 854534-10-4 CAPLUS  
 CN 4-Piperidinecarboxamide, 1-[1-(4-chlorophenyl)-2-(3-ethyl-1,2-dihydro-2-oxo-6-quinoxalinyloxy)ethyl]- (CA INDEX NAME)

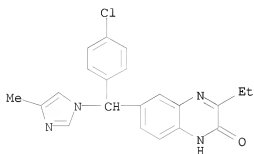


RN 854534-11-5 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[2-(4-chlorophenyl)-2-(4-hydroxy-4-phenyl-1-piperidinyl)ethyl]-3-ethyl- (CA INDEX NAME)



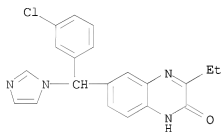
RN 854534-12-6 CAPLUS

CN 2-(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-(4-methyl-1H-imidazol-1-yl)methyl]-3-ethyl- (CA INDEX NAME)



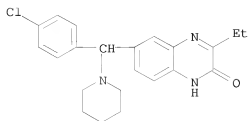
RN 854534-13-7 CAPLUS

CN 2-(1H)-Quinoxalinone, 6-[(3-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-ethyl- (CA INDEX NAME)



RN 854534-14-8 CAPLUS

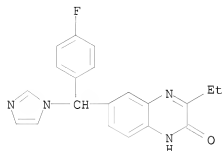
CN 2-(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1-piperidinylmethyl]-3-ethyl- (CA INDEX NAME)





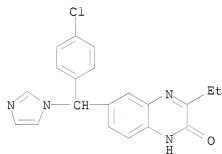
RN 854534-15-9 CAPLUS

CN 2(1H)-Quinoxalinone, 3-ethyl-6-[(4-fluorophenyl)-1H-imidazol-1-ylmethyl]-  
(CA INDEX NAME)



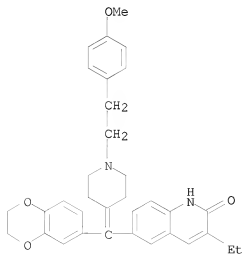
RN 854534-16-0 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-ethyl-  
(CA INDEX NAME)

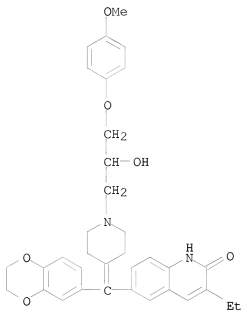


RN 854534-20-6 CAPLUS

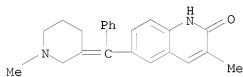
CN 2(1H)-Quinolone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl) [1-[2-(4-methoxyphenyl)ethyl]-4-piperidinylidene]methyl]-3-ethyl- (CA INDEX NAME)



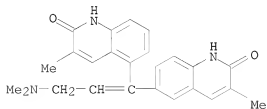
RN 854534-21-7 CAPLUS  
 CN 2(1H)-Quinolinsonone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl) [1-[2-hydroxy-3-(4-methoxyphenoxy)propyl]-4-piperidinylidene)methyl]-3-ethyl- (CA INDEX NAME)



RN 854534-22-8 CAPLUS  
 CN 2(1H)-Quinolinsonone, 3-methyl-6-[(1-methyl-3-piperidinylidene)phenylmethyl]- (CA INDEX NAME)

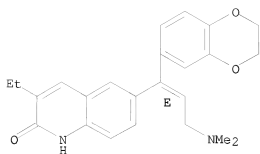


RN 854534-23-9 CAPLUS  
 CN 2(1H)-Quinolinsonone, 5-[1-(1,2-dihydro-3-methyl-2-oxo-6-quinoliny)-3-(dimethylamino)-1-propen-1-yl]-3-methyl- (CA INDEX NAME)



RN 854534-24-0 CAPLUS  
 CN 2(1H)-Quinolinsonone, 6-[(1E)-1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(dimethylamino)-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

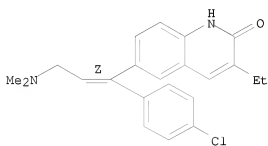
Double bond geometry as shown.



RN 854534-25-1 CAPLUS

CN 2(1H)-Quinolinone, 6-[(1Z)-1-(4-chlorophenyl)-3-(dimethylamino)-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

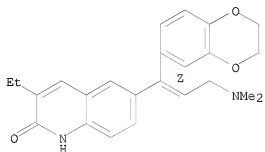
Double bond geometry as shown.



RN 854534-26-2 CAPLUS

CN 2(1H)-Quinolinone, 6-[(1Z)-1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(dimethylamino)-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

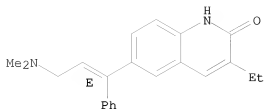
Double bond geometry as shown.



RN 854534-27-3 CAPLUS

CN 2(1H)-Quinolinone, 6-[(1E)-3-(dimethylamino)-1-phenyl-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

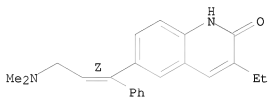
Double bond geometry as shown.



RN 854534-28-4 CAPLUS

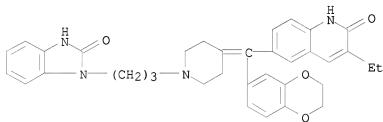
CN 2(1H)-Quinolinone, 6-[(1Z)-3-(dimethylamino)-1-phenyl-1-propen-1-yl]-3-ethyl- (CA INDEX NAME)

Double bond geometry as shown.



RN 854534-29-5 CAPLUS

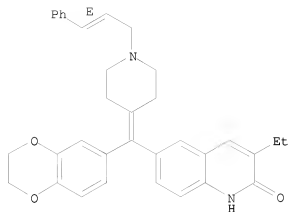
CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[1-[3-(2,3-dihydro-2-oxo-1H-benzimidazol-1-yl)propyl]-4-piperidinylidene]methyl]-3-ethyl- (CA INDEX NAME)



RN 854534-30-8 CAPLUS

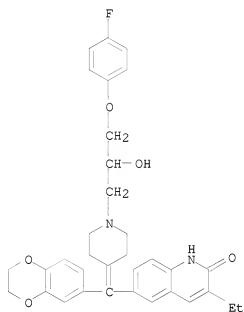
CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl)[1-[(2E)-3-phenyl-2-propen-1-yl]-4-piperidinylidene]methyl]-3-ethyl- (CA INDEX NAME)

Double bond geometry as shown.



RN 854534-31-9 CAPLUS

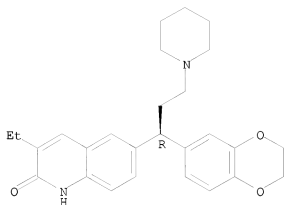
CN 2(1H)-Quinolinone, 6-[(2,3-dihydro-1,4-benzodioxin-6-yl) 1-[3-(4-fluorophenoxy)-2-hydroxypropyl]-4-piperidinylidene]methyl]-3-ethyl- (CA INDEX NAME)



RN 854534-32-0 CAPLUS

CN 2(1H)-Quinolinone, 6-[(1R)-1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(1-piperidinyl)propyl]-3-ethyl- (CA INDEX NAME)

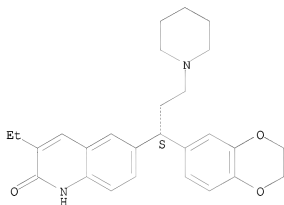
Absolute stereochemistry.



RN 854534-33-1 CAPLUS

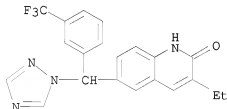
CN 2(1H)-Quinolinone, 6-[(1S)-1-(2,3-dihydro-1,4-benzodioxin-6-yl)-3-(1-piperidinyl)propyl]-3-ethyl- (CA INDEX NAME)

Absolute stereochemistry.



RN 854535-35-6 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[1H-1,2,4-triazol-1-yl[3-(trifluoromethyl)phenyl]methyl]- (CA INDEX NAME)



IT 854534-37-5P 854534-38-6P 854534-40-0P

854534-42-2P 854534-48-8P 854534-49-9P

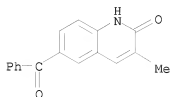
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854534-53-5P 854534-62-6P 854534-64-8P

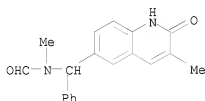
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of 6-alkenyl and 6-phenylalkyl substituted 2-quinolinones and 2-quinoxalinones as poly(ADP-ribose) polymerase inhibitors)

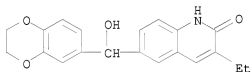
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 CN 2(1H)-Quinolinone, 6-benzoyl-3-methyl- (CA INDEX NAME)



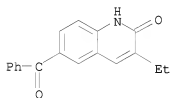
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 CN Formamide, N-[(1,2-dihydro-3-methyl-2-oxo-6-quinolinyl)phenylmethyl]-N-methyl- (CA INDEX NAME)



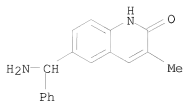
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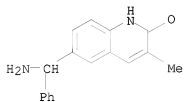
RN 854534-42-2 CAPLUS  
 CN 2(1H)-Quinolinone, 6-benzoyl-3-ethyl- (CA INDEX NAME)



RN 854534-48-8 CAPLUS  
 CN 2(1H)-Quinolinone, 6-(aminophenylmethyl)-3-methyl- (CA INDEX NAME)

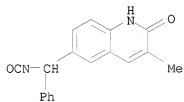


RN 854534-49-9 CAPLUS  
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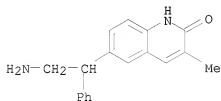


● HCl

RN 854534-50-2 CAPLUS  
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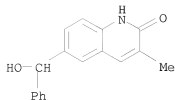


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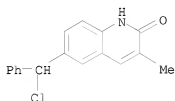


RN 854534-52-4 CAPLUS  
 CN 2 (1H)-Quinolinone, 6-(hydroxyphenylmethyl)-3-methyl- (CA INDEX NAME)

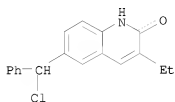




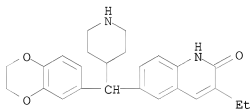
RN 854534-53-5 CAPLUS  
 CN 2(1H)-Quinolinone, 6-(chlorophenylmethyl)-3-methyl- (CA INDEX NAME)



RN 854534-62-6 CAPLUS  
 CN 2(1H)-Quinolinone, 6-(chlorophenylmethyl)-3-ethyl- (CA INDEX NAME)



RN 854534-64-8 CAPLUS  
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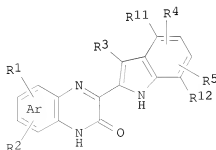
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L7 ANSWER 3 OF 15 CAPLUS COPYRIGHT 2010 ACS ON STN  
 ACCESSION NUMBER: 2004:430796 CAPLUS  
 DOCUMENT NUMBER: 141:7139  
 TITLE: Preparation of indolylquinoxalinones for treating hyperproliferative disorders and diseases associated with angiogenesis

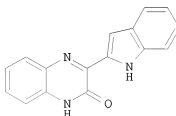
INVENTOR(S): Ladouceur, Gaetan H.; Bear, Brian; Bi, Cheng;  
 Brittelli, David R.; Burke, Michael J.; Chen, Gang;  
 Cook, James; Dumas, Jacques; Sibley, Robert; Turner,  
 Michael R.  
 PATENT ASSIGNEE(S): Bayer Pharmaceuticals Corporation, USA  
 SOURCE: PCT Int. Appl., 217 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004043950	A1	20040527	WO 2003-US36003	20031110 <--
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RW:	BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
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BR 2003016169	A	20050927	BR 2003-16169	20031110 <--
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MX 2005004779	A	20050722	MX 2005-4779	20050504 <--
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PRIORITY APPLN. INFO.:			US 2002-425490P	P 20021112 <--
			US 2003-460915P	P 20030407 <--
			US 2003-484202P	P 20030630 <--
			WO 2003-US36003	W 20031110 <--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT  
 OTHER SOURCE(S): MARPAT 141:7139  
 GI



I



II

AB The invention relates to title compds. I [wherein Ar = 6-membered aromatic ring containing 0-2 N atoms; R1 and R2 = independently H, halo, CF3, acyl, piperidinyl, piperazinyl, morpholinyl, or (un)substituted alkyl, alkoxy,

amino, pyrrolidinyl, Ph, etc.; R3 = H, alkyl, OH, NO2, NH2, alkylamino, alkoxyamino, or (un)substituted benzoylamino; R4 = H, OH, halo, CN, acyl, sulfamoyl, trialkylsiloxy, tetrazolyl, thienyl, pyrrolyl, pyrimidinyl, oxazolyl, furanyl, or (un)substituted alkyl, alkenyl, alkynyl, alkoxy, amino, oxadiazolyl, Ph, pyridyl(oxy), carbamoyl; R11 and R12 = independently H, F, or Cl with the proviso that when one of R11 and R12 = F or Cl, the other must be H; and pharmaceutically acceptable salts and esters thereof]. The invention also relates to the use of I and their pharmaceutical compns. for treating hyperproliferative disorders and diseases associated with angiogenesis (no data). Examples include representative syntheses for compds. of the invention, pharmaceutical compns. comprising them, and tumor model assays (no specific data given). For instance, N-Boc-indole was coupled with di-Me oxalate using t-BuLi to give tert-Bu 2-[methoxy(oxo)acetyl]-1H-indole-1-carboxylate (72%). Cyclization of the dione with 1,2-phenylenediamine in AcOH afforded the quinoxalinone II (77%).

IT 694531-84-5P 694531-85-6P 694531-86-7P  
694531-93-6P 694532-04-2P

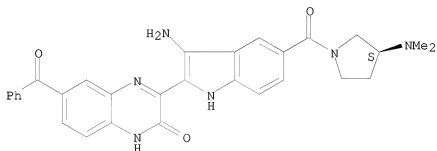
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(antiproliferative and angiogenesis inhibitor; preparation of indolylquinoxalinones for treating hyperproliferative disorders and diseases associated with angiogenesis)

RN 694531-84-5 CAPLUS

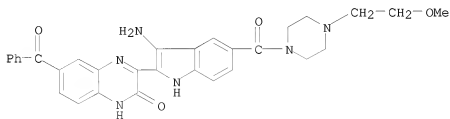
CN 2(1H)-Quinoxalinone, 3-[3-amino-5-[[ (3S)-3-(dimethylamino)-1-pyrrolidinyl]carbonyl]-1H-indol-2-yl]-6-benzoyl- (CA INDEX NAME)

Absolute stereochemistry.



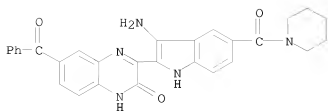
RN 694531-85-6 CAPLUS

CN 2(1H)-Quinoxalinone, 3-[3-amino-5-[[4-(2-methoxyethyl)-1-piperazinyl]carbonyl]-1H-indol-2-yl]-6-benzoyl- (CA INDEX NAME)



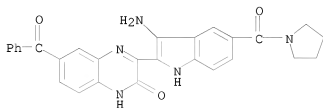
RN 694531-86-7 CAPLUS

CN 2(1H)-Quinoxalinone, 3-[3-amino-5-(1-piperidinylcarbonyl)-1H-indol-2-yl]-6-benzoyl- (CA INDEX NAME)



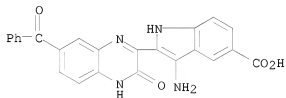
RN 694531-93-6 CAPLUS

CN 2(1H)-Quinoxalinone, 3-[3-amino-5-(1-pyrrolidinylcarbonyl)-1H-indol-2-yl]-6-benzoyl- (CA INDEX NAME)



RN 694532-04-2 CAPLUS

CN 1H-Indole-5-carboxylic acid, 3-amino-2-(7-benzoyl-3,4-dihydro-3-oxo-2-quinoxaliny)- (CA INDEX NAME)



OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD (9 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 4 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

ACCESSION NUMBER: 2003:796538 CAPLUS

DOCUMENT NUMBER: 139:323440

TITLE: Preparation of radiolabeled quinolines and quinolinones as metabotropic glutamate receptor mGluR1 antagonists for use in positron emission tomography.

INVENTOR(S): Lesage, Anne Simone Josephine; Bischoff, Francois Paul; Janssen, Cornelius Gerardus Maria; Lavreysen, Hilde

PATENT ASSIGNEE(S): Janssen Pharmaceutica N.V., Belg.

SOURCE: PCT Int. Appl., 148 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

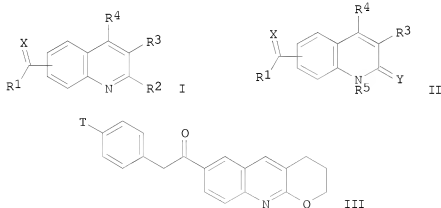
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003082350	A2	20031009	WO 2003-EP3240	20030326 <--
WO 2003082350	A3	20040304		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GO, GW, ML, MR, NE, SN, TD, TG				
CA 2479109	A1	20031009	CA 2003-2479109	20030326 <--
AU 2003226737	A1	20031013	AU 2003-226737	20030326 <--
AU 2003226737	B2	20080904		
BR 2003008945	A	20050104	BR 2003-8945	20030326 <--
EP 1492571	A2	20050105	EP 2003-745282	20030326 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
CN 1642580	A	20050720	CN 2003-807387	20030326 <--
JP 2005524679	T	20050818	JP 2003-579882	20030326 <--
NZ 535438	A	20060831	NZ 2003-535438	20030326 <--
IN 2004DN02631	A	20050401	IN 2004-DN2631	20040908 <--
US 20060083676	A1	20060420	US 2004-509069	20040924 <--
US 7517517	B2	20090414		
MX 2004009435	A	20050125	MX 2004-9435	20040928 <--
ZA 2004007820	A	20051011	ZA 2004-7820	20040928 <--
NO 2004004635	A	20041027	NO 2004-4635	20041027 <--
PRIORITY APPLN. INFO.:				
			EP 2002-76254	A 20020329 <--
			WO 2003-EP3240	W 20030326 <--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT  
OTHER SOURCE(S): MARPAT 139:323440  
GI



AB Radiolabeled title compds. [I, II; X = O, S, C(R6)2, NR7; Y = O, S; R1 = (substituted) alkyl, cycloalkyl, cycloalkylalkyl, thienyl, quinolinyl, etc.; R2 = H, halo, cyano, alkyl, amino, heterocyclyl, etc.; R3, R4 = H, halo, OH, cyano, alkyl, alkoxy, etc.; R2R3 = (CH2)3-6, Z4CH2CH2CH2, Z4CH2CH2, etc.; Z4 = O, S, SO2, NR11; R11 = H, alkyl, PhCH2, alkoxy carbonyl; R3R4 = (CH2)4, CH:CHCH:CH; R5 = H, cycloalkyl, piperidinyl, oxothienyl, tetrahydrothienyl, aralkyl, alkoxyalkyl, etc.; R6

= H, aryl, alkyl, aminoalkyl; R7 = amino, OH], were prepared Most preferred are radiolabeled compds. in which the radioactive isotope is selected from <sup>3</sup>H, <sup>11</sup>C and <sup>18</sup>F. The invention also relates to their use in a diagnostic method, in particular for marking and identifying a mGluR1 receptor in biol. material, as well as to their use for imaging an organ, in particular using positron emission tomog. (PET). Thus, title compound (III) was prepared by tritiation of the corresponding bromide in THF using tritium gas and Pd/C catalyst. The purified product showed specific activity of 25 Ci/mmol.

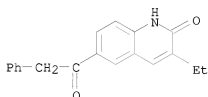
IT 409344-47-4P 409344-48-5P 409344-56-5P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of radiolabeled quinolines and quinolinones as metabotropic glutamate receptor mGluR1 antagonists for use in positron emission tomog.)

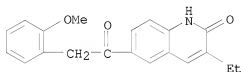
RN 409344-47-4 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-(2-phenylacetyl)- (CA INDEX NAME)



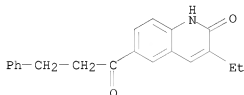
RN 409344-48-5 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[2-(2-methoxyphenyl)acetyl]- (CA INDEX NAME)



RN 409344-56-5 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-(1-oxo-3-phenylpropyl)- (CA INDEX NAME)



OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (4 CITINGS)

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 5 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

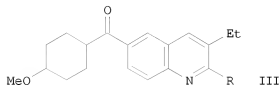
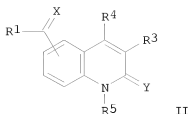
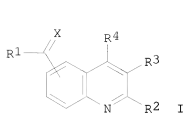
ACCESSION NUMBER: 2002:275968 CAPLUS

DOCUMENT NUMBER: 136:309857

TITLE: Preparation of quinolines and quinolinones as

INVENTOR(S): metabotropic glutamate receptor antagonists  
Mabire, Dominique Jean-Pierre; Venet, Marc Gaston;  
Coupa, Sophie; Poncelet, Alain Philippe; Lesage, Anne  
Simone Josephine  
PATENT ASSIGNEE(S): Janssen Pharmaceutica N.V., Belg.  
SOURCE: PCT Int. Appl., 114 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2002028837	A1	20020411	WO 2001-EP11135	20010925 <--
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
CA 2421782	A1	20020411	CA 2001-2421782	20010925 <--
AU 2001093847	A	20020415	AU 2001-93847	20010925 <--
BR 2001014253	A	20030701	BR 2001-14253	20010925 <--
EP 1332133	A1	20030806	EP 2001-974298	20010925 <--
EP 1332133	B1	20080709		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
HU 2003002167	A2	20031028	HU 2003-2167	20010925 <--
JP 2004510764	T	20040408	JP 2002-532423	20010925 <--
NZ 524945	A	20050128	NZ 2001-524945	20010925 <--
EE 200300126	A	20050415	EE 2003-126	20010925 <--
EE 5195	B1	20090817		
CN 1703403	A	20051130	CN 2001-816717	20010925 <--
AU 2001293847	B2	20070524	AU 2001-293847	20010925 <--
AT 400558	T	20080715	AT 2001-974298	20010925 <--
ES 2309095	T3	20081216	ES 2001-974298	20010925 <--
IL 155163	A	20090803	IL 2001-155163	20010925 <--
TW 306854	B	20090301	TW 2001-90124220	20011002 <--
KR 818965	B1	20080404	KR 2003-702014	20030211 <--
HR 2003000229	A2	20030630	HR 2003-229	20030324 <--
IN 2003MN00328	A	20050211	IN 2003-MN328	20030324 <--
BG 107672	A	20040130	BG 2003-107672	20030326 <--
ZA 2003002515	A	20040630	ZA 2003-2515	20030331 <--
NO 2003001474	A	20030505	NO 2003-1474	20030401 <--
NO 325079	B1	20080128		
MX 2003002907	A	20030624	MX 2003-2907	20030401 <--
US 20040082592	A1	20040429	US 2003-381987	20030814 <--
US 7115630	B2	20061003		
US 20050209273	A1	20050922	US 2005-133678	20050520 <--
US 7629468	B2	20091208		
PRIORITY APPLN. INFO.:			EP 2000-203419	A 20001002 <--
			WO 2001-EP11135	W 20010925 <--
			US 2003-381987	A3 20030814 <--
OTHER SOURCE(S):	MARPAT	136:309857		
GI				

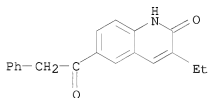


AB The title compds. [I or II; X = O, C(R6)2; (wherein R6 = H, aryl, alkyl, etc.); R1 = alkyl, aryl, thienyl, etc.; R2 = H, halo, CN, etc.; R3, R4 = H, alkyl; or R2 and R3 may be taken together to form (CH2)3, (CH2)4, CH:CHCH:CH, etc.; or R3 and R4 may be taken together to form CH:CHCH:CH, (CH2)4; R5 = H, cycloalkyl, piperidinyl, etc.; Y = O, S; or Y and R5 may be taken together to form CH:NN, N:NN, NCH:CH], useful for treating or preventing glutamate-induced diseases of the central nervous system, were prepared. Thus, reacting cis-III [R = Cl] with SnMe4 in the presence of Pg(PPh3)4 in PhMe afforded 17% cis-III [R = Me] which showed antagonism at a dose of 2.5 mg/kg bodyweight in cold allodynia test in rats with a Bennett ligation.

IT 409344-47-4P 409344-48-5P 409344-56-5P  
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
 (preparation of quinolines and quinolinones as metabotropic glutamate receptor antagonists)

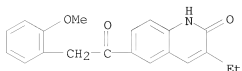
RN 409344-47-4 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-(2-phenylacetyl)- (CA INDEX NAME)



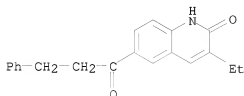
RN 409344-48-5 CAPLUS

CN 2(1H)-Quinolinone, 3-ethyl-6-[2-(2-methoxyphenyl)acetyl]- (CA INDEX NAME)





RN 409344-56-5 CAPLUS  
CN 2(1H)-Quinolinone, 3-ethyl-6-(1-oxo-3-phenylpropyl)- (CA INDEX NAME)



OS.CITING REF COUNT: 18 THERE ARE 18 CAPLUS RECORDS THAT CITE THIS  
RECORD (18 CITINGS)  
REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 6 OF 15 CAPLUS COPYRIGHT 2010 ACS ON STN  
ACCESSION NUMBER: 2001:476381 CAPLUS  
DOCUMENT NUMBER: 135:100164  
TITLE: (R)-1-{(2-Oxo-1,2-dihydroquinolin-6-yl)[3-(trifluoromethyl)phenyl]methyl}-1H-1,2,4-triazol-4-ium bromide  
AUTHOR(S): Peeters, Oswald M.; Blaton, Norbert M.; De Ranter, Camiel J.  
CORPORATE SOURCE: Faculteit Farmaceutische Wetenschappen, Laboratorium voor Analytische Chemie en Medicinale Fysicochemie, Katholieke Universiteit Leuven, Louvain, B-3000, Belg.  
SOURCE: Acta Crystallographica, Section E: Structure Reports Online (2001), E57(7), o655-o656  
CODEN: ACSEBH; ISSN: 1600-5368  
URL: <http://journals.iucr.org/e/issues/2001/07/00/ya6033/ya6033.pdf>  
PUBLISHER: International Union of Crystallography  
DOCUMENT TYPE: Journal; (online computer file)  
LANGUAGE: English

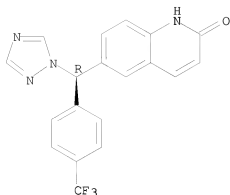
AB The metabolism of all-trans-retinoic acid is mediated by a cytochrome dependent P 450 system. The title compound, C19H14F3N4O+·Br- (R111214), is an inhibitor of P 450. The three planar ring systems, viz. the triazolyl, Ph and quinolinone groups, are arranged in a propeller-like fashion around the central CH group. The dihedral angles formed by the triazolyl/phenyl, triazolyl/quinolinone and phenyl/quinolinone planes are 55.8(1), 79.85(9) and 78.49(9)°, resp. The N-H...O H bonds, involving the triazolium N-H group and the quinolinone O atom, link the cations into infinite chains stretching along the c axis of the crystal. Crystallog. data are given.

IT 349553-99-7  
RL: PRP (Properties)  
(crystal structure of)

RN 349553-99-7 CAPLUS

CN 2(1H)-Quinolinone, 6-[(R)-1H-1,2,4-triazol-1-yl][4-(trifluoromethyl)phenyl]methyl]-, hydrobromide (1:1) (CA INDEX NAME)

Absolute stereochemistry.



● HBr

REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 7 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

ACCESSION NUMBER: 2001:223060 CAPLUS

DOCUMENT NUMBER: 135:5590

TITLE: Some nucleophilic reactions with 6-benzoyl-2,3-dichloroquinoxaline: synthesis of tetrazolo[1,5-a]quinoxaline, 2-methylidene-1,3-dithiolo[4,5-b]quinoxalines, quinoxalino[2,3-b]quinoxalines and pyrazolo[1',5':1,2]imidazolo[4,5-b]-quinoxalines

AUTHOR(S): El-Gaby, M. S. A.; El-Sharief, A. M. Sh; Ammar, Y. A.; Mohamed, Y. A.; El-Salam, A. A. Abd

CORPORATE SOURCE: Department of Chemistry, Faculty of Science, Al-Azhar University at Assiut, Assiut, 71524, Egypt

SOURCE: Indian Journal of Chemistry, Section B: Organic Chemistry Including Medicinal Chemistry (2001), 40B(3), 195-200

CODEN: IJSBDB; ISSN: 0376-4699

PUBLISHER: National Institute of Science Communication, CSIR Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 135:5590

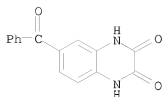
AB The starting material 6-benzoyl-2,3-dichloroquinoxaline is subjected to some nucleophilic reagents to study the effect of the benzoyl group on the reactivity of the two chlorine atoms.

IT 143702-68-5

RL: RCT (Reactant); RACT (Reactant or reagent) (reactions of 6-benzoyl-2,3-dichloroquinoxaline with nucleophiles)

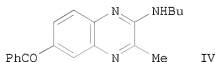
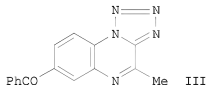
RN 143702-68-5 CAPLUS

CN 2,3-Quinoxalinedione, 6-benzoyl-1,4-dihydro- (CA INDEX NAME)



OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD  
(5 CITINGS)  
REFERENCE COUNT: 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

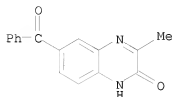
L7 ANSWER 8 OF 15 CAPLUS COPYRIGHT 2010 ACS ON STN  
ACCESSION NUMBER: 2000:527827 CAPLUS  
DOCUMENT NUMBER: 134:162992  
TITLE: Synthesis and antimicrobial activities of some novel  
quinoxalinone derivatives  
AUTHOR(S): Ali, M. M.; Ismail, M. M. F.; El-Gaby, M. S. A.;  
Zahran, M. A.; Ammar, Y. A.  
CORPORATE SOURCE: Dep. of Chemistry, Faculty of Science, Al-Azhar Univ.,  
Cairo, 11884, Egypt  
SOURCE: Molecules [online computer file] (2000),  
5(6), 864-873  
CODEN: MOLEFW; ISSN: 1420-3049  
URL: <http://www.mdpi.org/molecules/papers/50600864.pdf>  
PUBLISHER: Molecular Diversity Preservation International  
DOCUMENT TYPE: Journal; (online computer file)  
LANGUAGE: English  
OTHER SOURCE(S): CASREACT 134:162992  
GI



AB Condensation of 4-benzoyl-1,2-phenylenediamine with sodium pyruvate in acetic acid furnished two products, which were identified as 6-benzoyl- (I) and 7-benzoyl-3-methyl-2(1H)-quinoxalinone (II). Fusion of I with aromatic aldehydes furnished the styryl derivs. Alkylation of I and II with di-Me sulfate or Et chloroacetate produced the N-alkyl derivs. Hydrazinolysis of one ester derivative with hydrazine hydrate afforded the hydrazide derivative, which underwent condensation with aldehydes to give the corresponding hydrazone derivs. In addition, chlorination of I with thionyl chloride afforded the 2-chloro derivative, which was subjected to reaction with sodium azide and n-butylamine to yield the corresponding tetrazolo (III) and n-butylamino (IV) derivs., resp. The structures of the compds. prepared were confirmed by anal. and spectral data. Also, some of the synthesized compds. were screened for antimicrobial activity.

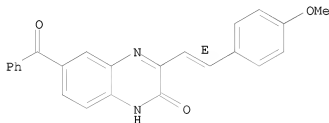
IT 325469-51-0P  
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and antimicrobial activities of quinoxalinone derivs.)

RN 325469-51-0 CAPLUS  
CN 2(1H)-Quinoxalinone, 6-benzoyl-3-methyl- (CA INDEX NAME)



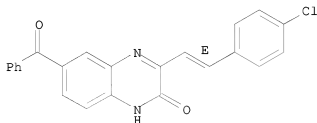
IT 325469-54-3P  
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)  
(preparation and antimicrobial activities of quinoxalinone derivs.)  
RN 325469-54-3 CAPLUS  
CN 2(1H)-Quinoxalinone, 6-benzoyl-3-[(1E)-2-(4-methoxyphenyl)ethenyl]- (CA INDEX NAME)

Double bond geometry as shown.



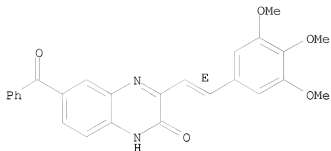
IT 325469-53-2P 325469-55-4P  
RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation and antimicrobial activities of quinoxalinone derivs.)  
RN 325469-53-2 CAPLUS  
CN 2(1H)-Quinoxalinone, 6-benzoyl-3-[(1E)-2-(4-chlorophenyl)ethenyl]- (CA INDEX NAME)

Double bond geometry as shown.



RN 325469-55-4 CAPLUS  
CN 2(1H)-Quinoxalinone, 6-benzoyl-3-[(1E)-2-(3,4,5-trimethoxyphenyl)ethenyl]- (CA INDEX NAME)

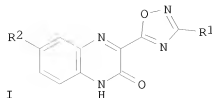
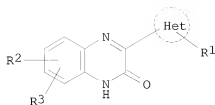
Double bond geometry as shown.



OS.CITING REF COUNT: 41 THERE ARE 41 CAPLUS RECORDS THAT CITE THIS  
RECORD (41 CITINGS)  
REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 9 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN  
ACCESSION NUMBER: 1998:672545 CAPLUS  
DOCUMENT NUMBER: 129:275932  
ORIGINAL REFERENCE NO.: 129:56265a,56268a  
TITLE: Preparation of 3-oxadiazolylquinoxaline derivatives  
having affinity to benzodiazepine receptor  
INVENTOR(S): Ohno, Kazunori; Odai, Osamu; Furukawa, Kiyoshi; Oka,  
Makoto  
PATENT ASSIGNEE(S): Dainippon Pharmaceutical Co., Ltd., Japan  
SOURCE: PCT Int. Appl., 43 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: Japanese  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9842701	A1	19981001	WO 1998-JP827	19980227 <--
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE,				
DK, EE, ES, FI, GB, GE, GH, GM, GW, HU, ID, IL, IS, JP, KE, KG,				
KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO,				
NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA,				
UG, US, UZ, VN, YU, ZW				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI,				
FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM,				
GA, GN, ML, MR, NE, SN, TD, TG				
JP 2002241379	A	20020828	JP 1997-87646	19970321 <--
AU 9861179	A	19981020	AU 1998-61179	19980227 <--
PRIORITY APPLN. INFO.:				
			JP 1997-87646	A 19970321 <--
			WO 1998-JP827	W 19980227 <--
OTHER SOURCE(S): MARPAT 129:275932				
GI				



II

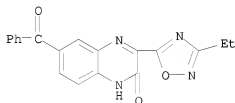
AB Novel 3-oxadiazolylquinoxaline derivs. represented by general formula (I; wherein Het is oxadiazolyl; R1 is hydrogen, lower alkyl, trifluoromethyl, lower cycloalkyl, lower alkenyl, lower alkynyl, optionally substituted aryl, optionally substituted heteroaryl, or lower alkoxy; R2 is hydrogen, lower alkyl, trifluoromethyl, lower cycloalkyl, halogeno, hydroxy, lower alkoxy, cyano, nitro, acyl, optionally substituted benzoyl, amino, lower mono- or dialkylamino, lower alkoxycarbonylmethoxy, lower mono- or dialkylaminocarbonylmethoxy, or optionally substituted benzyloxy; and R3 is hydrogen, lower alkyl, lower cycloalkyl, halogeno, or lower alkoxy), which are useful as a medicine, in particular, which have a selective affinity for benzodiazepine receptors and are useful as a brain activator and a remedy for senile dementia and Alzheimer's disease. Thus, a solution of 1,2-dihydro-2-oxo-3-quinoxalinecarboxylic acid and N,N'-carbonyl diimidazole in DMF was heated with stirring for 3 h at 60°, followed by adding acetamidoxime, and the stirring was continued for another 1.5 h to give 52.6% the title compound (II; R1 = Me; R2 = H). The latter compound and I (R1 = Et, R2 = OMe) inhibited the binding of [3H]diazepam to synaptosome membrane fraction prepared from rat brain with IC50 of 11.5 and 1.41 nM, resp.

IT 213743-73-8P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of oxadiazolylquinoxaline derivs. having affinity to benzodiazepine receptor as brain activators and remedies for senile dementia and Alzheimer's disease)

RN 213743-73-8 CAPLUS

CN 2(1H)-Quinoxalinone, 6-benzoyl-3-(3-ethyl-1,2,4-oxadiazol-5-yl)- (CA INDEX NAME)



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)  
REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 10 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

ACCESSION NUMBER: 1998:210752 CAPLUS

DOCUMENT NUMBER: 128:257445

ORIGINAL REFERENCE NO.: 128:50967a,50970a

TITLE: Preparation of indolylbenzoquinoxalinones and related compounds as protein kinase C inhibitors.

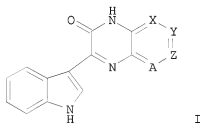
INVENTOR(S): Bergstrand, Hakan; Karabelas, Kostas; Sjo, Peter  
 PATENT ASSIGNEE(S): Astra Aktiebolag (Publ), Swed.  
 SOURCE: PCT Int. Appl., 63 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9813368	A1	19980402	WO 1997-SE1582	19970919 <--
W:	AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW			
RW:	GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG			
IN 1997DE02638	A	20050311	IN 1997-DE2638	19970916 <--
TW 472045	B	20020111	TW 1997-86113549	19970918 <--
ZA 9708469	A	19980325	ZA 1997-8469	19970919 <--
CA 2265854	A1	19980402	CA 1997-2265854	19970919 <--
AU 9744775	A	19980417	AU 1997-44775	19970919 <--
AU 716279	B2	20000224		
EP 929551	A1	19990721	EP 1997-943259	19970919 <--
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO			
NZ 334531	A	20000929	NZ 1997-334531	19970919 <--
US 6271231	B1	20010807	US 1997-981266	19971218 <--
US 20010025043	A1	20010927	US 2001-865231	20010525 <--
PRIORITY APPLN. INFO.:			SE 1996-3505	A 19960925 <--
			SE 1997-2747	A 19970718 <--
			WO 1997-SE1582	A 19970919 <--
			US 1997-981266	A3 19971218 <--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

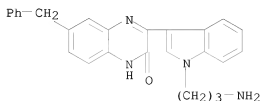
OTHER SOURCE(S): MARPAT 128:257445

GI



AB Title compds. [I; A, X, Y, Z = C, N;  $\geq 2$  of A, X, Y, Z = C; may be substituted and/or annulated; excluding 3-(1H-indol-3-yl)-1H-quinoxalin-2-one, 3-(2-methyl-1H-indol-3-yl)-1H-quinoxalin-2-one, and 3-(1,2-diphenyl-1H-indol-3-yl)-1H-quinoxalin-2-one], were prepared as protein kinase C inhibitors (no data). Thus, 1,2-phenylenediamine was stirred overnight with [1-[3-(1,3-dioxoisindol-2-yl)propyl]-1H-indol-3-yl]oxoacetic acid 2,5-dioxopyrrolidin-1-yl ester (preparation given) in THF to give 3-[3-(3-oxo-3,4-dihydroquinoxalin-2-yl)indol-1-yl]propylammonium

acetate. The latter was stirred with MeNH<sub>2</sub> in THF/H<sub>2</sub>O to give  
 3-[3-(3-oxo-3,4-dihydroquinoxalin-2-yl)indol-1-yl]propylammonium acetate.  
 IT 205377-77-1P  
 RL: BAC (Biological activity or effector, except adverse); BSU (Biological  
 study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);  
 BIOL (Biological study); PREP (Preparation); USES (Uses)  
 (preparation of indolylbenzoquinoxalinones and related compds. as protein  
 kinase C inhibitors)  
 RN 205377-77-1 CAPLUS  
 CN 2(1H)-Quinoxalinone, 3-[1-(3-aminopropyl)-1H-indol-3-yl]-6-(phenylmethyl)-  
 , 2,2,2-trifluoroacetate (1:1) (CA INDEX NAME)  
 CM 1  
 CRN 205377-76-0  
 CMF C26 H24 N4 O



CM 2  
 CRN 76-05-1  
 CMF C2 H F3 O2



OS.CITING REF COUNT: 10 THERE ARE 10 CAPLUS RECORDS THAT CITE THIS  
 RECORD (11 CITINGS)  
 REFERENCE COUNT: 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS  
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 11 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN  
 ACCESSION NUMBER: 1996:527663 CAPLUS  
 DOCUMENT NUMBER: 125:167994  
 ORIGINAL REFERENCE NO.: 125:31485a,31488a  
 TITLE: Preparation of  
 6-[triazolyl(3-trifluoromethylphenyl)methyl]-2-  
 quinolin(thi)ones for treatment of keratinization  
 disorders  
 INVENTOR(S): Venet, Marc Gaston; Mabire, Dominique Jean-Pierre;  
 Sanz, Gerard Charles  
 PATENT ASSIGNEE(S): Janssen Pharmaceutica N.V., Belg.  
 SOURCE: PCT Int. Appl., 32 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1



PATENT INFORMATION:

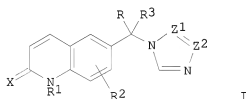
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9620200	A1	19960704	WO 1995-EP5173	19951221 <--
W: AL, AM, AU, BB, BG, BR, BY, CA, CN, CZ, EE, FI, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LS, LT, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, RO, RU, SD, SG, SI, SK, TJ, TM, TT, UA, UG, US, UZ, VN				
RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
IN 1995CA01685	A	20050304	IN 1995-CA1685	19951220 <--
CA 2207268	A1	19960704	CA 1995-2207268	19951221 <--
AU 9644362	A	19960719	AU 1996-44362	19951221 <--
AU 698199	B2	19981029		
EP 800524	A1	19971015	EP 1995-943237	19951221 <--
EP 800524	B1	20011031		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE				
CN 1171789	A	19980128	CN 1995-197162	19951221 <--
CN 1085668	C	20020529		
JP 10511654	T	19981110	JP 1995-520222	19951221 <--
BR 9510504	A	19990601	BR 1995-10504	19951221 <--
RU 2165419	C2	20010420	RU 1997-112898	19951221 <--
AT 207924	T	20011115	AT 1995-943237	19951221 <--
PT 800524	E	20020429	PT 1995-943237	19951221 <--
ES 2166838	T3	20020501	ES 1995-943237	19951221 <--
PL 182956	B1	20020531	PL 1995-321041	19951221 <--
ZA 9510989	A	19970627	ZA 1995-10989	19951227 <--
IL 116577	A	20000229	IL 1995-116577	19951227 <--
US 5922734	A	19990713	US 1997-860239	19970616 <--
FI 9702794	A	19970627	FI 1997-2794	19970627 <--
NO 9703029	A	19970627	NO 1997-3029	19970627 <--
NO 311220	B1	20011029		

PRIORITY APPLN. INFO.:

EP 1994-203773	A	19941228 <--
WO 1995-EP5173	W	19951221 <--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 125:167994  
GI



AB Title compds. [I; R = 3-(F3C)C6H4][II; R1 = H, NH2, alkyl; R2,R3 = H, halo, alkyl; X = O or S; 1 of Z1,Z2 = N and the other = CH] were prepared. Thus, (R)-II (R1-R3 = H, X = O, Z1 = N, Z2 = CH) gave complete suppression of estradiol undecylate-induced vaginal keratinization in 50% of ovariectomized rats at 1.25mg/kg orally.

IT 180421-65-2P	180421-66-3P	180421-67-4P
180421-68-5P	180421-69-6P	180421-70-9P
180421-71-0P	180421-72-1P	180421-73-2P
180421-74-3P	180421-75-4P	

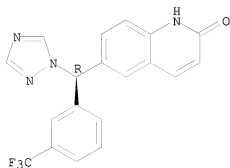
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of 6-[triazolyl(3-trifluoromethylphenyl)methyl]-2-quinolin(thi)ones for treatment of keratinization disorders)

RN 180421-65-2 CAPLUS

CN 2(1H)-Quinolinone, 6-[1H-1,2,4-triazol-1-yl[3-(trifluoromethyl)phenyl]methyl]-, (R)- (9CI) (CA INDEX NAME)

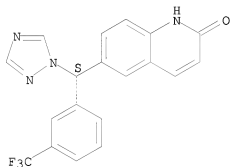
Absolute stereochemistry. Rotation (-).



RN 180421-66-3 CAPLUS

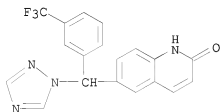
CN 2(1H)-Quinolinone, 6-[1H-1,2,4-triazol-1-yl[3-(trifluoromethyl)phenyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



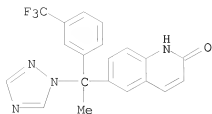
RN 180421-67-4 CAPLUS

CN 2(1H)-Quinolinone, 6-[1H-1,2,4-triazol-1-yl[3-(trifluoromethyl)phenyl]methyl]- (CA INDEX NAME)



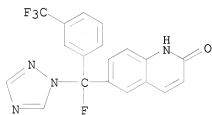
RN 180421-68-5 CAPLUS

CN 2(1H)-Quinolinone, 6-[1-[(1H-1,2,4-triazol-1-yl)-1-[3-(trifluoromethyl)phenyl]ethyl]- (CA INDEX NAME)



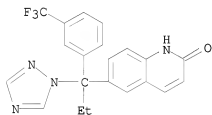
RN 180421-69-6 CAPLUS

CN 2(1H)-Quinolinone, 6-[fluoro-1H-1,2,4-triazol-1-yl]-3-(trifluoromethyl)phenylmethyl- (CA INDEX NAME)



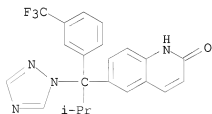
RN 180421-70-9 CAPLUS

CN 2(1H)-Quinolinone, 6-[1-(1H-1,2,4-triazol-1-yl)-1-[3-(trifluoromethyl)phenyl]propyl]- (CA INDEX NAME)



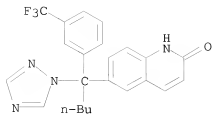
RN 180421-71-0 CAPLUS

CN 2(1H)-Quinolinone, 6-[2-methyl-1-(1H-1,2,4-triazol-1-yl)-1-[3-(trifluoromethyl)phenyl]propyl]- (CA INDEX NAME)

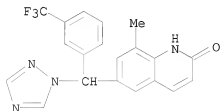


RN 180421-72-1 CAPLUS

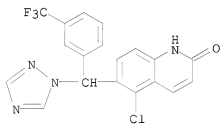
CN 2(1H)-Quinolinone, 6-[1-(1H-1,2,4-triazol-1-yl)-1-[3-(trifluoromethyl)phenyl]pentyl]- (CA INDEX NAME)



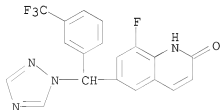
RN 180421-73-2 CAPLUS  
 CN 2(1H)-Quinolinone, 8-methyl-6-[1H-1,2,4-triazol-1-yl][3-(trifluoromethyl)phenyl]methyl- (CA INDEX NAME)



RN 180421-74-3 CAPLUS  
 CN 2(1H)-Quinolinone, 5-chloro-6-[1H-1,2,4-triazol-1-yl][3-(trifluoromethyl)phenyl]methyl- (CA INDEX NAME)



RN 180421-75-4 CAPLUS  
 CN 2(1H)-Quinolinone, 8-fluoro-6-[1H-1,2,4-triazol-1-yl][3-(trifluoromethyl)phenyl]methyl- (CA INDEX NAME)

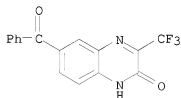


OS.CITING REF COUNT: 3 THERE ARE 3 CAPLUS RECORDS THAT CITE THIS RECORD (3 CITINGS)  
 REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 12 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN  
 ACCESSION NUMBER: 1992:592207 CAPLUS  
 DOCUMENT NUMBER: 117:192207  
 ORIGINAL REFERENCE NO.: 117:33223a,33226a  
 TITLE: Fluorine-19 NMR studies on the mechanism of riboflavin  
 synthase. Synthesis of  
 6-(trifluoromethyl)-7-oxo-8-(D-ribityl)lumazine and  
 6-(trifluoromethyl)-7-methyl-8-(D-ribityl)lumazine  
 AUTHOR(S): Cushman, Mark; Patel, Hemantkumar H.; Scheuring,  
 Johannes; Bacher, Adelbert  
 CORPORATE SOURCE: Sch. Pharm. Pharm. Sci., Purdue Univ., West Lafayette,  
 IN, 47907, USA  
 SOURCE: Journal of Organic Chemistry (1992), 57(21),  
 5630-43  
 CODEN: JOCEAH; ISSN: 0022-3263  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 GI

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

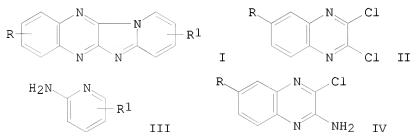
AB Title oxo-(D-ribityl)lumazine I was synthesized by reaction of Me  
 trifluoropyruvate with 5-amino-6-(D-ribitylamino)pyrimidine-2,4(1H,3H)-  
 dione hydrochloride and utilized as a 19F NMR probe of the light  
 riboflavin synthase of Bacillus subtilis. I was found to be an inhibitor  
 of riboflavin synthase with an inhibition constant KI = 55 µM. The  
 enzyme-bound ligand gave rise to several broad 19F NMR signals which were  
 shifted to low field. The bound ligand I could be displaced from the  
 enzyme by the enzyme product, riboflavin (II), and the product analog,  
 5-nitroso-6-(ribitylamino)-2,4(1H,3H)-pyrimidinedione. Title  
 methyl-(D-ribityl)lumazine III was synthesized by reaction of  
 5-amino-6-(D-ribitylamino)pyrimidine-2,4(1H,3H)-dione hydrochloride with  
 1,1,1-trifluorobutane-2,3-dione. Three mols. of III can be bound  
 relatively tightly per mol of riboflavin synthase, i.e., one ligand mol.  
 per protein subunit. A scheme for the catalytic cycle of riboflavin  
 synthase is proposed.  
 IT 143309-79-9P  
 RL: SPN (Synthetic preparation); PREP (Preparation)  
 (preparation of)  
 RN 143309-79-9 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-benzoyl-3-(trifluoromethyl)- (CA INDEX NAME)



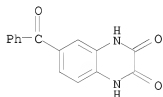
OS.CITING REF COUNT: 33 THERE ARE 33 CAPLUS RECORDS THAT CITE THIS  
 RECORD (33 CITINGS)

L7 ANSWER 13 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN  
 ACCESSION NUMBER: 1992:571381 CAPLUS  
 DOCUMENT NUMBER: 117:171381  
 ORIGINAL REFERENCE NO.: 117:29633a,29636a

TITLE: Synthesis of pyrido[1',2':1,2]imidazo[4,5-b]quinoxalines  
 AUTHOR(S): Tanaka, Kiyoshi; Takahashi, Hideki; Takimoto, Kozo; Sugita, Masahiko; Mitsuhashi, Keiyo  
 CORPORATE SOURCE: Fac. Eng., Seikei Univ., Musahino, 180, Japan  
 SOURCE: Journal of Heterocyclic Chemistry (1992), 29(4), 771-7  
 CODEN: JHTCAD; ISSN: 0022-152X  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 OTHER SOURCE(S): CASREACT 117:171381  
 GI



AB Synthesis of title compds. I (R = H, 8-, 9-Cl, 8-, 9-Bz, 8-, 9-NO<sub>2</sub>; R<sub>1</sub> = H, 1-, 2-, 3-, 4-Me, 4-PhCH<sub>2</sub>O) by the facile cyclizations of 2,3-dichloroquinoxalines II with 2-aminopyridines III and of 2-amino-3-chloroquinoxalines IV (R ≠ H) with various substituted pyridines is described.  
 IT 143702-68-5  
 RL: RCT (Reactant); RACT (Reactant or reagent) (chlorination of)  
 RN 143702-68-5 CAPLUS  
 CN 2,3-Quinoxalinedione, 6-benzoyl-1,4-dihydro- (CA INDEX NAME)



OS.CITING REF COUNT: 15 THERE ARE 15 CAPLUS RECORDS THAT CITE THIS RECORD (16 CITINGS)

L7 ANSWER 14 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN  
 ACCESSION NUMBER: 1990:612014 CAPLUS  
 DOCUMENT NUMBER: 113:212014  
 ORIGINAL REFERENCE NO.: 113:35835a, 35838a  
 TITLE: Preparation of (1H-azol-1-ylmethyl)quinolines, -quinazolines, and -quinoxalines as drugs  
 INVENTOR(S): Freyne, Eddy Jean Edgard; Venet, Marc Gaston; Raeymaekers, Alfons Herman Margaretha; Sanz, Gerard Charles  
 PATENT ASSIGNEE(S): Janssen Pharmaceutica N. V., Belg.  
 SOURCE: Eur. Pat. Appl., 106 pp.  
 CODEN: EPXXDW

DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 371564	A2	19900606	EP 1989-203014	19891128 <--
EP 371564	A3	19910529		
EP 371564	B1	19950712		
R: AT, BE, CH, DE, ES, FR, GB, GR, IT, LI, LU, NL, SE				
US 5028606	A	19910702	US 1989-434957	19891113 <--
US 5037829	A	19910806	US 1989-435120	19891113 <--
CA 2002864	A1	19900529	CA 1989-2002864	19891114 <--
CA 2002864	C	19991116		
DK 8905994	A	19900530	DK 1989-5994	19891128 <--
DK 172748	B1	19990628		
NO 8904734	A	19900530	NO 1989-4734	19891128 <--
NO 174509	B	19940207		
NO 174509	C	19940518		
AU 8945646	A	19900607	AU 1989-45646	19891128 <--
AU 620946	B2	19920227		
HU 52498	A2	19900728	HU 1989-6220	19891128 <--
HU 205106	B	19920330		
ZA 8909076	A	19910731	ZA 1989-9076	19891128 <--
SU 1780536	A3	19921207	SU 1989-4742543	19891128 <--
IL 92486	A	19930708	IL 1989-92486	19891128 <--
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CN 1042912	A	19900613	CN 1989-108925	19891129 <--
CN 1033752	C	19970108		
JP 02223579	A	19900905	JP 1989-307793	19891129 <--
JP 2916181	B2	19990705		
US 5151421	A	19920929	US 1991-672298	19910320 <--
US 5185346	A	19930209	US 1991-704746	19910523 <--
US 5268380	A	19931207	US 1992-973871	19921110 <--
US 5441954	A	19950815	US 1993-131817	19931005 <--
CN 1106004	A	19950802	CN 1994-117801	19941102 <--
CN 1036002	C	19971001		
CN 1106005	A	19950802	CN 1994-117802	19941102 <--
CN 1036003	C	19971001		
US 5612354	A	19970318	US 1995-409551	19950323 <--

PRIORITY APPLN. INFO.:

GB 1988-27820	A	19881129 <--
GB 1988-27821	A	19881129 <--
GB 1988-27822	A	19881129 <--
US 1989-434205	B2	19891113 <--
US 1989-434957	A3	19891113 <--
US 1991-704746	A3	19910523 <--
US 1992-973871	A3	19921110 <--
US 1993-131817	A3	19931005 <--

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 113:212014

GI For diagram(s), see printed CA Issue.

AB The title compds. [I; R = H, alkyl; X1:X2 = CH:CH, CH:N, N:CH; Y = H, alkyl, cycloalkyl, alkenyl, alkynyl, (un)substituted aryl, aralkyl; Z = (un)substituted (oxo)quinolinyl, (oxo- or thioxo)quinazolinyl, (oxo- or dioxo)quinoxalinyl] were prepared as retinoic acid metabolism inhibitors, aromatase inhibitors, etc. Thus, 3,4-dihydroquinolin-2(1H)-one was stirred 2 h at 70° with BzCl in DMF containing AlCl3 and the product reduced by NaBH4 to give hydroxymethylquinolinone II (R1 = Ph, R2 = OH). II (R1 = Me, R2 = OH) was stirred overnight with SOCl2 in THF and the

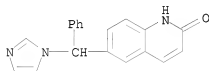
product II (R1 = Me, R2 = Cl) stirred overnight at 60-70° with 1H-imidazole in DMSO to give II (R1 = Me, R2 = imidazolo) which maintained plasma levels of i.v. administered all-trans-retinoic acid at  $\geq 10$  ng/mL in rats 2 h after oral administration of 40 mg/kg.

IT	120067-41-6P	130344-00-2P	130344-01-3P
	130344-02-4P	130344-03-5P	130346-18-8P
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	130347-44-3P	130347-45-4P	130347-46-5P
	130347-47-6P	130347-48-7P	130347-62-5P

RL: SPN (Synthetic preparation); PREP (Preparation)  
(preparation of, as retinoate metabolism and aromatase inhibitor)

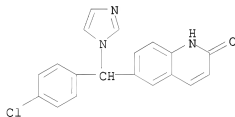
RN 120067-41-6 CAPLUS

CN 2(1H)-Quinolinone, 6-(1H-imidazol-1-ylphenylmethyl)- (CA INDEX NAME)



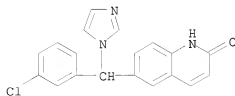
RN 130344-00-2 CAPLUS

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RN 130344-01-3 CAPLUS

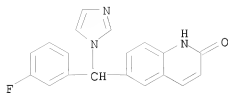
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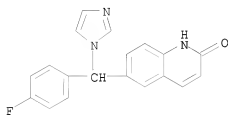
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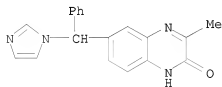




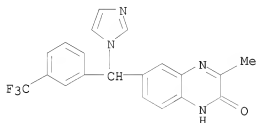
RN 130344-03-5 CAPLUS  
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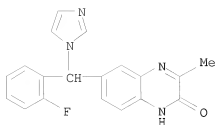
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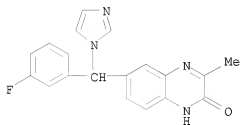
RN 130346-22-4 CAPLUS  
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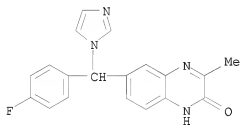
RN 130346-25-7 CAPLUS  
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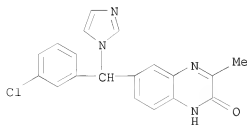
RN 130346-26-8 CAPLUS  
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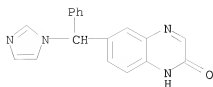
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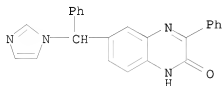
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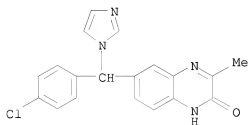
RN 130346-42-8 CAPLUS  
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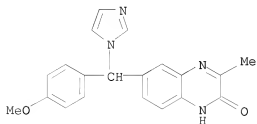
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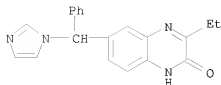
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 CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-methyl- (CA INDEX NAME)



RN 130346-68-8 CAPLUS  
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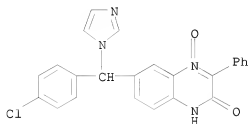


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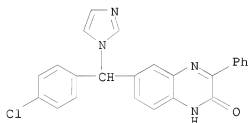
RN 130347-21-6 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-phenyl-, 4-oxide (CA INDEX NAME)



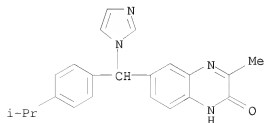
RN 130347-22-7 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-phenyl-, (CA INDEX NAME)



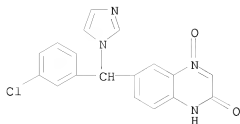
RN 130347-23-8 CAPLUS

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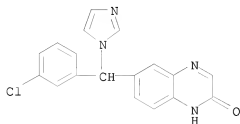


RN 130347-25-0 CAPLUS

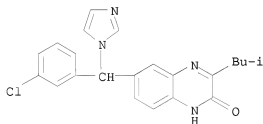
CN 2(1H)-Quinoxalinone, 6-[(3-chlorophenyl)-1H-imidazol-1-ylmethyl]-, 4-oxide (CA INDEX NAME)



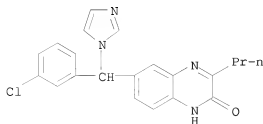
RN 130347-26-1 CAPLUS  
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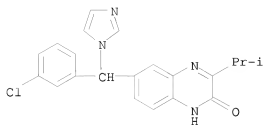
RN 130347-28-3 CAPLUS  
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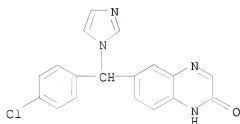
RN 130347-30-7 CAPLUS  
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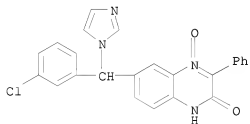
RN 130347-31-8 CAPLUS  
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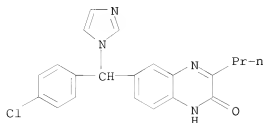
RN 130347-33-0 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]- (CA INDEX NAME)



RN 130347-35-2 CAPLUS  
 CN 2(1H)-Quinoxalinone, 6-[(3-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-phenyl-, 4-oxide (CA INDEX NAME)

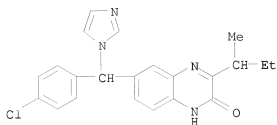


RN 130347-37-4 CAPLUS  
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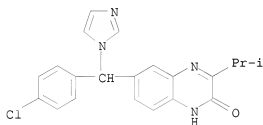
RN 130347-39-6 CAPLUS

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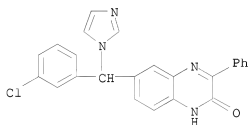
RN 130347-41-0 CAPLUS

CN 2(1H)-Quinoxalinone, 6-[(4-chlorophenyl)-1H-imidazol-1-ylmethyl]-3-(1-methylethyl)- (CA INDEX NAME)



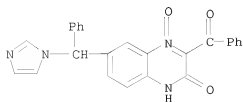
RN 130347-42-1 CAPLUS

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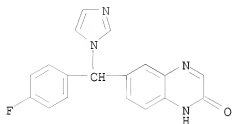
RN 130347-44-3 CAPLUS

CN 2(1H)-Quinoxalinone, 3-benzoyl-6-(1H-imidazol-1-ylphenylmethyl)-, 4-oxide (CA INDEX NAME)

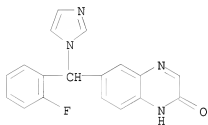


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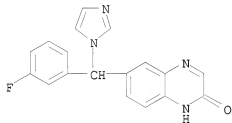
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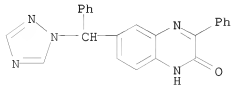
RN 130347-46-5 CAPLUS  
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RN 130347-47-6 CAPLUS  
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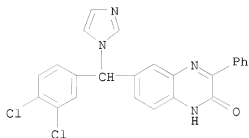
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CN 2(1H)-Quinoxalinone, 3-phenyl-6-(phenyl-1H-1,2,4-triazol-1-ylmethyl)- (CA INDEX NAME)



RN 130347-62-5 CAPLUS  
CN 2(1H)-Quinoxalinone, 6-[(3,4-dichlorophenyl)-1H-imidazol-1-ylmethyl]-3-



phenyl- (CA INDEX NAME)



OS.CITING REF COUNT: 24 THERE ARE 24 CAPLUS RECORDS THAT CITE THIS RECORD (43 CITINGS)

L7 ANSWER 15 OF 15 CAPLUS COPYRIGHT 2010 ACS on STN

ACCESSION NUMBER: 1989:407401 CAPLUS

DOCUMENT NUMBER: 111:7401

ORIGINAL REFERENCE NO.: 111:1422h,1423a

TITLE: Imidazole- or pyridine-containing carbostyryls as combined thromboxane synthetase and cyclic-AMP phosphodiesterase inhibitors, their preparation, and pharmaceuticals containing them  
Walker, Keith A. M.; Bruno, John J.; Martinez, Gregory R.

INVENTOR(S): Syntex (U.S.A.), Inc., USA

PATENT ASSIGNEE(S): U.S., 20 pp.

SOURCE: CODEN: USXXAM

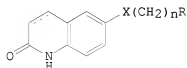
DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 4792561	A	19881220	US 1986-868845	19860529 <--
US 4921862	A	19900501	US 1988-247134	19880921 <--
PRIORITY APPLN. INFO.:			US 1986-868845	A3 19860529 <--
ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT				
OTHER SOURCE(S): CASREACT 111:7401; MARPAT 111:7401				
GI				



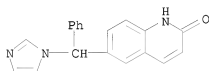
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AB Title compds. I [X = R1CR2, cis- or trans-CR3:CR4; R1 = H when R2 = OH, or R1 = Ph, phenylalkyl when R2 = H, OH; Ph is optionally monosubstituted; or R1R2 = O, C1-6 alkylidene, (substituted) benzylidene; R3 = H, C1-6 alkyl; R4 = H; R3R4 = bond; n = 0-3; R = 1-imidazolyl; dotted line = optional covalent bond] are prepared as thromboxane synthetase and cAMP phosphodiesterase inhibitors for treatment of disease characterized by elevated thromboxane levels or an imbalance of prostacyclin/thromboxane levels (no data). A mixture of CuI 11.6, (Ph3P)2PdCl2 86,

N-propargylimidazole (preparation given) 774 mg, and 6-bromo-3,4-dihydrocarbostyryl 1.5 g was stirred in 10mL pyridine and 2 mL triethylamine at 100° for 48 h under N. The reaction mixture was then treated with saturated aqueous K2CO3, extracted with 10% MeOH in CH2Cl2, and worked up to give 6-[3-(imidazol-1-yl)-1-propyn-1-yl]-3,4-dihydrocarbostyryl. The latter (502 mg) was stirred under H in the presence of 200 mg 10% Pd/C to give 6-[3-(imidazol-1-yl)propyl]-3,4-dihydrocarbostyryl (II). A tablet was formulated containing II 25, cornstarch 20, spray-dried lactose 153, and Mg stearate 2 mg.

IT 120067-41-6P  
 RL: SPN (Synthetic preparation); PREP (Preparation)  
 (preparation of, as cAMP phosphodiesterase and thromboxane synthetase inhibitors)

RN 120067-41-6 CAPLUS  
 CN 2(1H)-Quinolinone, 6-(1H-imidazol-1-ylphenylmethyl)- (CA INDEX NAME)



OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD (6 CITINGS)  
 REFERENCE COUNT: 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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FULL ESTIMATED COST	92.77	478.03
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-12.75	-12.75

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LOGOFF? (Y)/N/HOLD:y

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